

Page 1
Received: 12/12/86

SPECTRIX LOGO

REPORT

Work Order # 86-12-05/

01/22/87 12:47:29

REPORT Keystone Env. Resources, Inc.
TO Water Quality Engineering
440 College Park Drive
Monroeville, PA 15146-1536
ATTEN Shannon Craig

PREPARED Spectrix CorporationBY 3911 FondrenSuite 100Houston, Texas 77063-5821

ATTEN Sample Control
PHONE (713) 266-6800

H. Hernandez

CERTIFIED BY

CONTACT CRUZ

CLIENT KEYSTONE SAMPLES 10
COMPANY Keystone Company, Inc.
FACILITY _____

Please call the above number if you have any questions.

NOTE: ALL SAMPLES WILL BE RETAINED FOR 90 DAYS AND THEN
DISCARDED. IF YOU WISH YOUR SAMPLES RETURNED TO YOUR FACILITY
CALL SAMPLE CONTROL AT THE ABOVE NUMBER.

WORK ID Water
TAKEN Client
TRANS Client
TYPE Project 1573-73
P. O. # 14-5-50106
INVOICE under separate cover

SAMPLE IDENTIFICATION

- 01 LAB BLANK
02 CAV-DW02-001
03 CAV-DW08-001
04 CAV-DW09-001
05 CAV-DW14-001
06 SCK-MW04-001
07 SCK-MW10-001
08 SCK-MW13-001
09 SCK-MW22-001
10 TRIP BLANK

TEST CODES and NAMES used on this report

AG	Silver
AS	Argentic
BE	Berulium
CD	Cadmium
CN SU	Cuanide in Water
CR	Chromium
CU	Copper
FE	Iron
HG	Mercury
NI	Nickel
PB	Lead
PEST W	EPA Pesticides/PCB's-Water
SB	Antimony
SE	Selenium
TL	Thallium
ZN	Zinc

LABORATORY: SPECTRIX-HOUSTON
LABORATORY NO: 86-12-057
CLIENT: SPECTRIX-MONROEVILLE
SITE: SOUTH CAVALCADE

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	Inorganic Matrix Duplicate and Matrix Spike

004008

CHECKLIST

SECTION I - Support Information and QC Summaries

- Narrative
- Introduction to the Spectrix Organic Data Package
- Introduction to the Spectrix Inorganic Data Package
- Shipping Document, Chain-of-Custody
- ~~N/A~~ Organic MS/MSD Data Summary Form

004009

SPECTRIX
NARRATIVE - ORGANIC ANALYSES

CLIENT: SPECTRIX - MONROEVILLE
LAB NO: 8612057

SITE: South Cavalcade

-004010

I. RECEIPT

- A. DATE: 12/12/86
- B. NO. OF SAMPLES: 9
- C. SAMPLE TYPE: WATER
- D. ANALYSIS REQUESTED: ATT organics following EPA CLP protocol
- E. SHIPPING PROBLEMS: NONE
- F. DOCUMENTATION PROBLEMS: NONE
- G. QC DISTRIBUTION:

LEVEL	MATRIX	VOA	SV	PEST
LOW	WATER			
	SOIL			
MEDIUM	WATER			
	SOIL			

II. PREPARATION

- A. ANALYTICAL LEVELS: Low level

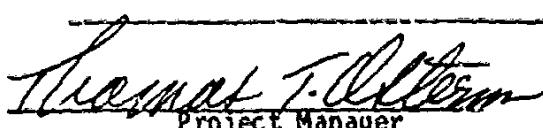
- B. EXTRACTION PROBLEMS: None

III. ANALYSIS

- A. ANALYSIS PROBLEMS: Some samples were diluted due to the high concentration of organics.

- B. SURROGATE RECOVERIES: Some surrogates were diluted out. The samples without dilution had surrogate recoveries within the range of method requirement.

- IV. OTHER PROBLEMS None


Thomas J. Oltman
Project Manager

SPECTRIX-HOUSTON

NARRATIVE - INORGANIC ANALYSES

CLIENT: Spectrix-Monroeville
LAB NO: 86-12-057

SITE: South Cavalcade

I. RECEIPT

- A. DATE: 12/12/86
B. NO. OF SAMPLES: Nine
C. SAMPLE TYPE: Water
D. ANALYSIS REQUESTED: Metals: Aq, As, Be, Cd, Cr, Cu, Fe, Hg, Ni, Pb, Sb, Se, Tl, Zn and CN

II. PREPARATION

A. REFERENCE METHOD: EPA CLP Protocol (SOW784)

B. ANALYTICAL LEVELS: Low level

C. DIGESTION PROBLEMS: None

III. ANALYSIS

A. REFERENCE METHOD: EPA CLP Protocol (SOW784)

B. ANALYSIS PROBLEMS: None

C. SPIKE RECOVERIES: None

IV. OTHER PROBLEMS None

Ken U. Easlo
Department Supervisor

004011

INTRODUCTION TO THE SPECTRIX ORGANIC DATA PACKAGE

The purpose of this introduction is to clarify several technical report sheets contained in this report. Data sheets showing analytes and concentration are called Organic Analysis Data Sheets (OADS). OADS pages 1 and 2 contain Hazardous Substance List (HSL) compounds associated with Volatile Organics and Semi-Volatile Organics (Base/Neutral/Acids). OADS pages 4 and 5 contain Tentatively Identified Compounds (TIC) found in the Volatiles and Semi-Volatiles analyses. All concentrations listed on pages 1 and 2 are stated in UG/L or UG/KG. Header information listed on pages 1 and 2 includes client number, Spectrix internal sample number, dilution factor used, instrument used in the analysis, date received/date analyzed and other pertinent information. Soil OADS also contain listings for percent moisture and soil pH determined during the extraction process. The reporting section for pages 1 and 2 of the OADS lists individual HSL compounds and four possible listings of the concentrations found. These four possibilities are the actual concentration found, nothing detected at the detection limit as noted with a "U", a value detected but just below the detection limit noted with a "J" and a value also found in the blank noted with a "B". It is possible that a sample may contain a value just below the detection limit and also found in the blank. In this instance a "JB" would be seen after a listed detection limit. A "J" value reported indicates a trace level was found below a concentration where the quantitated value may be in question.

OADS pages 4 and 5 in general list compounds not on the HSL but also found in the GC/MS analysis. These compounds or "non-targeted compounds" are identified by comparing the peak spectra in question to a list of several thousands library spectra stored on the GC/MS magnetic disc. A measure to which the library spectra and the sample spectra are similar is the "purity". Purity is listed on the OADS as a number between 1 and 1000 where 1000 would be an extremely confident identification. OADS pages 4 and 5 list two sections: a header and data reporting section. The header contains client sample name, instrument ID, analyst and data file name. The data section shows cas #, compound name, scan #, purity of each compound and UG/L or UG/KG found. The cas #, or chemical abstract service number, is an identifying number related to the organic compound name and can help clarify differences between common names and IUPAC terminology. The scan # relates to the location of the compound in the run (i.e., retention time). The concentration listed is estimated based on the Internal Standard concentrated which is added to the sample prior to analysis. These estimated values are listed as a "J" and are not to be confused with "J" listed on OADS pages 1 and 2. Only purity values greater than 750 will show the name and cas #. Purity values less than 750 are considered unknowns and only a scan # and concentration are listed. Unknown compounds on the OADS pages 4 and 5 will also be listed on individual library search spectra sheets included in the data package. These raw data sheets show the top 4 matches to the unknown compounds so that best estimation of the compound's identity can be ascertained by you the data user. Additional manual identifications may be achieved at additional costs.

Quantitation report sheets are also found in this package. These data sheets depict raw header and analysis data important for record keeping purposes. The quantitation report also lists surrogate recoveries showing the % recovery of similar HSL compounds for Quality Control determinations. EPA % recovery acceptance levels are available upon request.

INTRODUCTION TO THE SPECTRIX INORGANIC DATA PACKAGE

The inorganic portion of the data package is designed to serve two purposes, namely, explain the reporting format and describe the methods used in analysis of samples.

Inorganic Analytical Report

Metals are reported in micrograms per liter (ug/L) for aqueous samples and milligrams per kilogram dry weight (mg/Kg dry wt.) for non-aqueous samples. Non-aqueous samples have been calculated to their dry weight basis. Default units for each parameter are shown just below the individual test code. In situations where the reported units differs from the default unit, the reported units will be shown next to the reported value.

Example: The default units for BE_S is mg/Kg dry wt., which indicates that report will be in dry weight basis. Report that shows mg/Kg below a numerical value means that sample was not calculated to dry weight basis.

Percent solid of non-aqueous samples, if performed, will be reported as PSOLID in %.

Reporting for non-metallic elements follow same pattern as metals; ug/L or mg/L for aqueous samples and mg/Kg for non-aqueous samples. Here too, mg/Kg dry weight indicates that samples have been corrected for dry weight.

Methodology for Analysis of Inorganic Parameters

The analysis of metals in aqueous samples is performed according to EPA Contract Laboratory Program (CLP) Protocol Statement of Work (SOW) No. 785, "Inorganic Analysis of Multi-Media, Multi-Concentration Samples", July, 1985. Silver, beryllium, cadmium, chromium, copper, iron, nickel, and zinc are analyzed by ICP emission spectrometry. Antimony, arsenic, lead, selenium and thallium are determined by furnace - AA, and mercury is determined by manual cold vapor. All furnace - AA and ICP analyses are background corrected. L'vov platformed graphite tubes are used for all furnace - AA analyses and matrix modified where applicable.

The analysis of non-aqueous samples is done according to EPA CLP Protocol SOW No. 785. This is a modified version of methods in EPA manual "Test methods for Evaluating Solid Waste", EPA SW-846, July, 1982.

The EP Toxicity leaching procedure is performed according to methods published by EPA in "Test Methods for Evaluating Solid Wastes", EPA SW-846, July, 1982. Analysis of the leachates for metals is performed by the method for aqueous samples cited above.

004014

CHAIN OF CUSTODY RECORD

Client/Project Name Koppers Co Inc			Project Location South Cavalcade Houston		ANALYSES									
Project No. 1573-73			Field Logbook No. Scott Barnes / KER #2											
Sampler (Signature) Scott Barnes David (Hail)			Chain of Custody Tape No. NA											
Sample No./Identification	Date	Time	Lab Sample Number	Type of Sample	RN	AC	Lam	Plastic	Plastic	PCB	Test.	REMARKS		
CAV-0W09-001	86/12/12	1200		Ground Water	2	1	1	1						
SCK-MW04-001	86/12/12	1217		" "	2	1	1	1						
SCK-MW22-001	86/12/12	1430		" "	2	1	1	1						
Relinquished by: (Signature) Scott Barnes				Date 86/12/12	Time 1643	Received by: (Signature)				Date	Time			
Relinquished by: (Signature)				Date	Time	Received by: (Signature)				Date	Time			
Relinquished by: (Signature)				Date	Time	Received for Laboratory: (Signature) GMF Cruz				Date 12/2/84	Time 16:43			
Sample Disposal Method:				Disposed of by: (Signature) Monroeville, PA 15146								Date	Time	
SAMPLE COLLECTOR KER Environmental Research and Technology, Inc. 696 Virginia Road Concord, MA 01742 617-369-8910				ANALYTICAL LABORATORY Spectrix Corp. 3911 Fondren Suite 100 Houston, TX 77063								ERT		
												No. 7358		

004014

S104015

CHAIN OF CUSTODY RECORD

Client/Project Name <i>Koppers Co Inc</i>		Project Location <i>South Cascadade, Houston</i>		ANALYSES													
Project No. <i>1573-73</i>		Field Logbook No. <i>Scott Barnes KER #2</i>															
Sampler: (Signature) <i>Scott Barnes</i>		Chain of Custody Tape No. <i>NA</i>															
Sample No./Identification	Date	Time	Lab Sample Number	Type of Sample	BN	AW	Amber	Plastic	PC	PP	PC	Plastic	Amber	REMARKS			
SCK-MW16-001	86/12/12	0910		Ground water	2	1	1	1									
SCK-MW13-001	86/12/12	1400		" "	2	1	1	1									
CAP-00002-001	86/12/12	1330		" "	2	1	1	1									
Relinquished by: (Signature) <i>Scott Barnes</i>				Date <i>86/12/12</i>	Time <i>1644</i>	Received by: (Signature)				Date <i>86/12/12</i>	Time <i>1644</i>						
Relinquished by: (Signature)				Date	Time	Received by: (Signature)				Date	Time						
Relinquished by: (Signature)				Date	Time	Received for Laboratory: (Signature) <i>MF Cruz</i>				Date <i>86/12/12</i>	Time <i>1644</i>						
Sample Disposal Method:				Disposed of by: (Signature) <i>MF Cruz</i>								Date	Time				
SAMPLE COLLECTOR <i>KER</i> Environmental Research and Technology Inc. 696 Virginia Road Concord, MA 01742 617-369-8910 Monroeville, PA 15146				ANALYTICAL LABORATORY <i>Spectrix Corp</i> 3911 Fondren Suite 100 Houston, TX 77063								ERT					
												No. 7360					

910700

CHAIN OF CUSTODY RECORD

Client/Project Name <i>Koppers Co Inc</i>		Project Location <i>South Cavalcade, Houston</i>		ANALYSES							
Project No. <i>1573-73</i>		Field Logbook No. <i>Scott Barnes KER #2</i>		<input checked="" type="checkbox"/> Benzene	<input checked="" type="checkbox"/> <i>PCP</i>	<input checked="" type="checkbox"/> <i>PCB</i>	<input checked="" type="checkbox"/> <i>Lead</i>	<input checked="" type="checkbox"/> <i>Mercury</i>	<input checked="" type="checkbox"/> <i>Pesticides</i>	<input checked="" type="checkbox"/> <i>Plastics</i>	<input checked="" type="checkbox"/> <i>Rubber</i>
Sampler: (Signature) <i>Scott Barnes</i>		Chain of Custody Tape No. <i>NA</i>		REMARKS							
Sample No./Identification	Date	Time	Lab Sample Number	Type of Sample	<i>YOA samples preserved in the field 24 hours VDT 2 Trip Blanks</i>						
CAV-0W08-001	86/2/2	13105		Ground Water	2	1	1	1	1	2	
CAV-0W14-001	86/2/2	12402		" "	2	1	1	1	1	2	
CAV-0W09-001	86/2/2	1200		" "						2	
SCK-MW04-001	86/2/2	1217		" "						2	
SCK-MW22-001	86/2/2	1430		" "						2	
CAV-0W02-001	86/2/2	1320		" "						2	
SCK-MW13-001	86/2/2	1420		" "						2	
SCK-MW10-001	86/2/2	0910		" "						2	
Relinquished by: (Signature) <i>Scott Barnes</i>				Date 86/2/2	Time 1642	Received by: (Signature)			Date	Time	
Relinquished by: (Signature)				Date	Time	Received by: (Signature)			Date	Time	
Relinquished by: (Signature)				Date	Time	Received for Laboratory: (Signature) <i>MF Cruz</i>			Date 12/2/86	Time 16:45	
Sample Disposal Method:				Disposed of by: (Signature) <i>MF Cruz</i>					Date	Time	
SAMPLE COLLECTOR <i>KER</i> Environmental Research and Technology, Inc. 696 Virginia Road Concord, MA 01742 617-369-8910 <i>Monroeville, PA 15146</i>				ANALYTICAL LABORATORY <i>Spectrix Corp</i> 3911 Fondren Suite 100 <i>Houston, TX 77063</i>					ERT No. 7361		

CHECKLIST

SECTION II - Sample Data Packets

Sample No: CAV-OW02-001

- Organic Analysis Data Sheets (OADS)
 - VOA
 - SV
 - Pest.
- Tentatively Identified Compounds
 - VOA
 - SV
- Raw Data
 - VOA chromatogram
 - VOA quantitation report (surrogates)
 - SV chromatogram
 - SV quantitation report (surrogates)
 - Pest. chromatogram and integrator print out
- Inorganic Analysis Data Sheets (IADS)

004017

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-0W02-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 801205702
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *ypm*

CASE NO.: -----
QC REPORT NO.: -----
CONTRACT NO.: -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/18/86
DATE ANALYZED: 12/18/86

DATAFILE: 4U12057V02
DILUTION FACTOR: 1:100

CAS #	COMPOUND	UG/L	00
C010	CHLOROMETHANE	1000 U	1
C015	BROMOMETHANE	1000 U	0
C020	VINYL CHLORIDE	1000 U	0
C025	CHLOROETHANE	1000 U	0
C030	METHYLENE CHLORIDE	500 U	0
C035	ACETONE	1000 U	0
C040	CARBON DISULFIDE	500 U	0
C045	1, 1-DICHLOROETHENE	500 U	0
C050	1, 1-DICHLOROETHANE	500 U	0
C055	TRANS-1, 2-DICHLOROETHENE	500 U	0
C060	CHLOROFORM	500 U	0
C065	1, 2-DICHLOROETHANE	500 U	0
C110	2-BUTANONE	1000 U	0
C115	1, 1, 1-TRICHLOROETHANE	500 U	0
C120	CARBON TETRACHLORIDE	500 U	0
C125	VINYL ACETATE	1000 U	0
C130	BROMODICHLOROMETHANE	500 U	0
C140	1, 2-DICHLOROPROPANE	500 U	0
C145	TRANS-1, 3-DICHLOROPROPENE	500 U	0
C150	TRICHLOROETHENE	500 U	0
C155	DIBROMOCHLOROMETHANE	500 U	0
C160	1, 1, 2-TRICHLOROETHANE	500 U	0
C165	BENZENE	500 U	0
C170	CIS-1, 3-DICHLOROPROPENE	500 U	0
C175	2-CHLOROETHYL VINYL ETHER	1000 U	0
C180	BROMOFORM	500 U	0
C210	2-HEXANONE	1000 U	0
C295	4-METHYL-2-PENTANONE	1000 U	0
C220	TETRACHLOROETHENE	500 U	0
C225	1, 1, 2, 2-TETRACHLOROETHANE	500 U	0
C230	TOLUENE	500 U	0
C235	CHLOROBENZENE	500 U	0
C240	ETHYLBENZENE	500 U	0
C245	STYRENE	500 U	0
C250	TOTAL XYLENES	500 U	0

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: CAV-DW02 -001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205702
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *M.J.T.*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86...

SEMOVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: ... 12/10/86...
 DATE ANALYZED: 01/13/87

DATAFILE: 4U12057C02A

DILUTION FACTOR: 100.0....

CAS #	COMPOUND	UG/L	004019
C315	PHENOL	1000 U	
C325	BIS(2-CHLOROETHYL)ETHER	1000 U	
C330	2-CHLOROPHENOL	1000 U	
C335	1, 3-DICHLOROBENZENE	1000 U	
C340	1, 4-DICHLOROBENZENE	1000 U	
C345	BENZYL ALCOHOL	1000 U	
C350	1, 2-DICHLOROBENZENE	1000 U	
C355	2-METHYLPHENOL	870 J	
C360	BIS(2-CHLOROISOPROPYL)ETHER	1000 U	
C365	4-METHYLPHENOL	990 J	
C370	N-NITROSODIPropylAMINE	1000 U	
C375	HEXACHLOROETHANE	1000 U	
C410	NITROBENZENE	1000 U	
C415	ISOPHORONE	1000 U	
C420	2-NITROPHENOL	1000 U	
C425	2, 4-DIMETHYLPHENOL	4500	
C430	BENZOIC ACID	5000 U	
C435	BIS(2-CHLOROETHOXY)METHANE	1000 U	
C440	2, 4-DICHLOROPHENOL	1000 U	
C445	1, 2, 4-TRICHLOROBENZENE	1000 U	
C450	NAPHTHALENE	20000	
C455	4-CHLOROANILINE	1000 U	
C460	HEXACHLOROBUTADIENE	1000 U	
C465	P-CHLORO-M-CRESOL	1000 U	
C470	2-METHYLNAPHTHALENE	1200	
C510	HEXACHLOROCYCLOPENTADIENE	1000 U	
C515	2, 4, 6-TRICHLOROPHENOL	1000 U	
C520	2, 4, 5-TRICHLOROPHENOL	5000 U	
C525	2-CHLORONAPHTHALENE	1000 U	
C530	2-NITROANILINE	5000 U	
C535	DIMETHYL PHTHALATE	1000 U	
C540	ACENAPHTHYLENE	1000 U	
C545	3-NITROANILINE	5000 U	
C550	ACENAPHTHENE	770 J	
C555	2, 4-DINITROPHENOL	5000 U	
C560	4-NITROPHENOL	5000 U	
C565	DIBENZOFURAN	450 J	
C570	2, 4-DINITROTOLUENE	1000 U	
C575	2, 6-DINITROTOLUENE	1000 U	

SPECTRIX DC # ---- B

SAMPLE NUMBER: CAV-QW02-001

EMI VOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C02A

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	1000 U
C585	4-CHLOROPHENYL PHENYL ETHER	1000 U
C590	FLUORENE	350 J
C595	4-NITROANILINE	3000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	5000 U
C615	N-NITROSODIPHENYLAMINE	5000 U
C625	4-BROMOPHENYL PHENYL ETHER	1000 U
C630	HEXA CHLOROBENZENE	1000 U
C635	PENTACHLOROPHENOL	1000 U
C640	PHENANTHRENE	5000 U
C645	ANTHRACENE	330 J
C650	DI-N-BUTYL PHTHALATE	1000 U
C655	FLUORANTHENE	1000 U
C715	PYRENE	1000 U
C720	BUTYL BENZYL PHTHALATE	1000 U
C725	3, 3'-DICHLOROBENZIDINE	1000 U
C730	BENZO(A)ANTHRACENE	2000 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	1000 U
C740	CHRYSENE	1000 U
C760	DI-N-OCTYL PHTHALATE	1000 U
C765	BENZO(B)FLUORANTHENE	1000 U
C770	BENZO(K)FLUORANTHENE	1000 U
C775	BENZO(A)PYRENE	1000 U
C780	INDENO(1, 2, 3-CD)PYRENE	1000 U
C785	DIBENZO(A, H)ANTHRACENE	1000 U
C790	BENZO(GHI)PERYLENE	1000 U
		1000 U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

004020

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-02A

DC No: KEYSTONE
SMO No: CAV-DW02-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low

GPC Cleanup Yes ✓
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

Date Extracted: 12/18/86

Date Analyzed: 1/6/87

Dilution Factor: 1

Percent Moisture(decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOL-1016	0.50 U
11104-28-2	AROCLOL-1221	0.50 U
11141-16-5	AROCLOL-1232	0.50 U
53469-21-5	AROCLOL-1242	0.50 U
12672-29-3	AROCLOL-1248	0.50 U
11097-69-1	AROCLOL-1254	1.0 U
11096-82-5	AROCLOL-1260	1.0 U

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst.

6000

2nd Column Inst.

6000

N/A

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

Form 1

004021

INST ID: 4000

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-0W02-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ----

QC REPORT NO.: ----

ANALYST: TTO

DATAFILE: 4U12057V01

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
-------	-------------------------	-------	--------	--------

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

004052

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: CAV-DW02 - 001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: QU

DATAFILE: 4U12057C02A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
271-89-6	BENZOFURAN			UG/M
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	512	852	540 J
95-13-6	1H-INDENE	565	962	620 Q
526-75-0	PHENOL, 2,3-DIMETHYL-	576	931	1100 J
498-81-7	CYCLOHEXANEMETHANOL, . ALPHA., . ALPHA., 4-TR	655	850	710 J
526-75-0	PHENOL, 2,3-DIMETHYL-	700	846	1300 J
95-15-8	BENZO[B]THIOPHENE	729	910	4100 J
527-60-6	PHENOL, 2,4,6-TRIMETHYL-	752	896	2700 J
99-89-8	PHENOL, 4-(1-METHYLETHYL)-	764	896	750 J
1515-95-3	BENZENE, 1-ETHYL-4-METHOXY-	791	856	1200 J
3855-26-3	PHENOL, 2-ETHYL-4-METHYL-	801	886	1100 J
527-60-6	PHENOL, 2,4,6-TRIMETHYL-	823	850	2100 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	831	888	710 J
2471-83-2	1H-INDENE, 1-ETHYLDENE-	842		730 J
6351-10-6	1H-INDEN-1-OL, 2,3-DIHYDRO-	875	773	1500 J
197-34-8	PHENOL, 3,5-DIETHYL-	881	771	1000 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	908	843	630 J
700-12-9	BENZENE, PENTAMETHYL-	946		730 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1002	775	590 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1103		2200 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1214		1900 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

004024

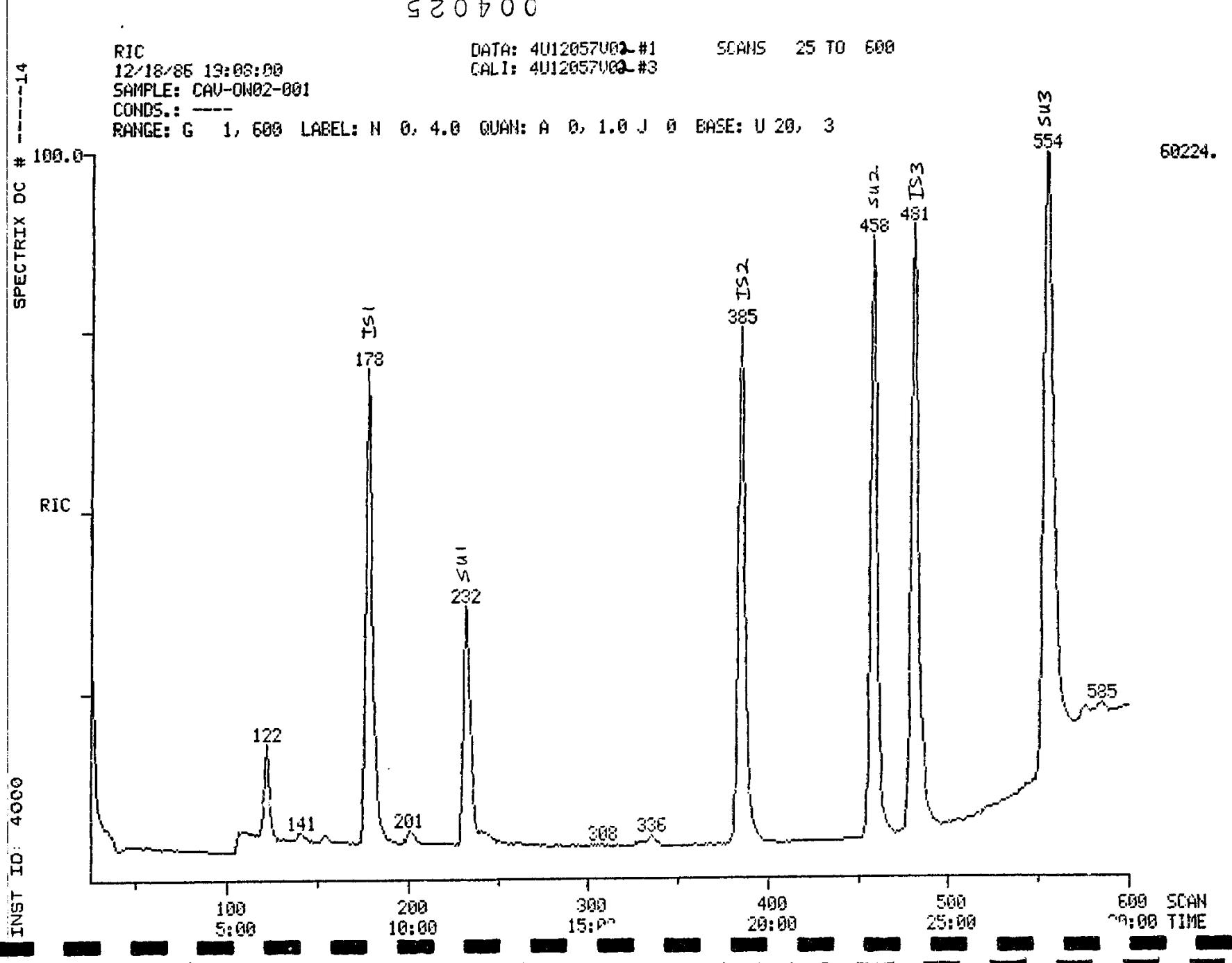
004025

RIC
12/18/86 19:00:00
SAMPLE: CAV-ON02-001
COND.: ---
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057U03-#1 SCANS 25 TO 600
CALI: 4U12057U03-#3

SPECTRIX DC # -----14

INST ID: 4000



004025

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID: CAV-OW02-001
FILENAME 4U12057VOA INST 4000
CLIENT KEYSTONE ANALYST TTO
DATE INJECTED 12/18/86 19:08:00 VERIFIED BY *NOT*
CALIB STD DATE 12-14-86 CORR. FACTOR 100.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA	
178 CI01	BROMOCHLOROMETHANE	**IS1**	128	5000 UG/L	1.000	16894
385 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	5000 UG/L	1.000	76118
481 CI20	D5-CHLOROBENZENE	**IS3**	117	5000 UG/L	1.000	65904
232 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	92 %	1.303	22214
458 CS05	DB-TOLUENE	**SU2**	98	97 %	0.952	67369
554 CS10	F-BROMOFLUOROBENZENE	**SU3**	95	104 %	1.152	50439
122 C030	METHYLENE CHLORIDE *		84	640 UG/L	0.685	4000

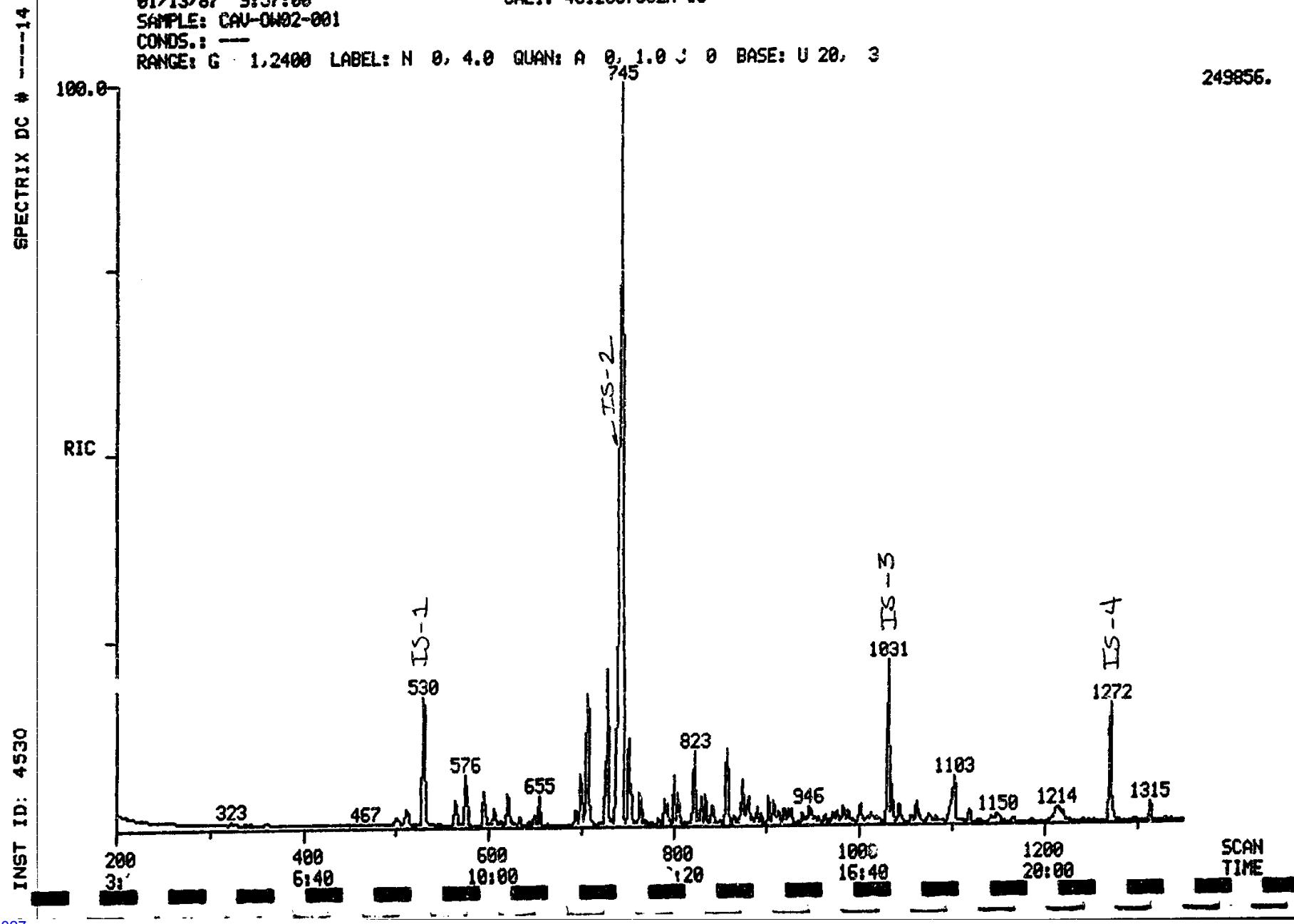
* BLANK CORRECTED ON DATA SHEET

004027

RIC
01/13/87 9:57:00
SAMPLE: CAV-0W02-001
COND.: —
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057C02A #1 SCANS 200 TO 1350
CALI: 4U12057C02A #3

249856.

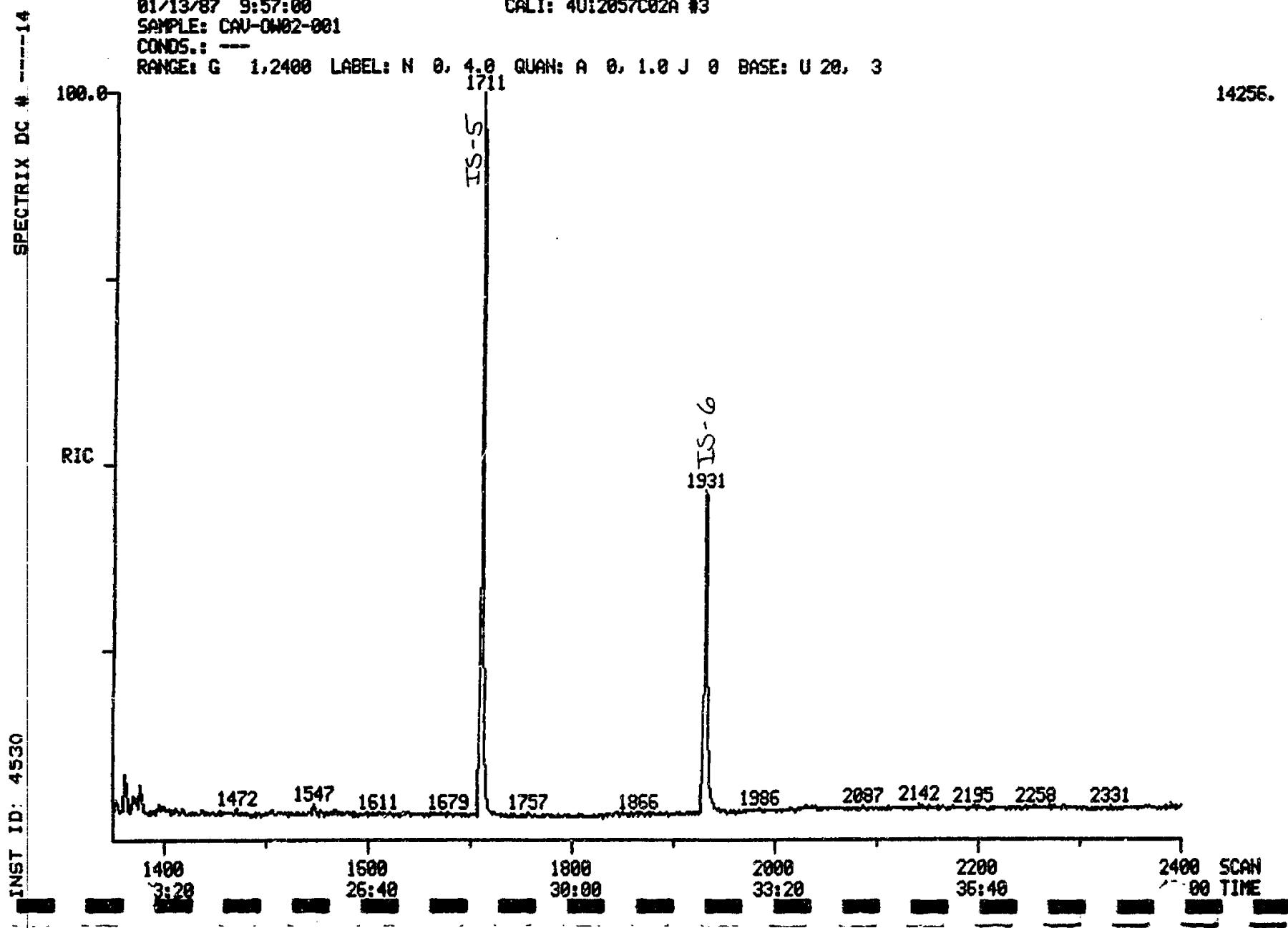


004028

RIC
01/13/87 9:57:00
SAMPLE: CAV-0W02-001
COND.: --
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057C02A #1 SCANS 1350 TO 2400
CALI: 4U12057C02A #3

14256.



004028

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	CAV-QW02-001		
FILENAME	4U12057C02A	INST ID.	4930
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/13/87 9:57:00	VERIFIED BY	<i>YPH</i>
CALIB STD DATE	01/09/87	CORR. FACTOR	100.00

SCANN#		SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
530	C130	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	4000 UG/L	1.000	21711
740	C140	NAPHTHALENE-D8 ** IS2 **	136	4000 UG/L	1.000	76447
1031	C150	ACENAPHTHENE-D10 ** IS3 **	164	4000 UG/L	1.000	38707
1272	C160	PHENANTHRENE-D10 ** IS4 **	188	4000 UG/L	1.000	45724
1711	C170	CHRYSENE-D12 ** IS5 **	240	4000 UG/L	1.000	21631
1931	C175	PERYLENE-D12 ** IS6 **	264	4000 UG/L	1.000	13444
395	C355	2-METHYLPHENOL	108	874 UG/L	1.123	7028
621	C365	4-METHYLPHENOL	108	989 UG/L	1.172	7840
707	C425	2,4-DIMETHYLPHENOL	122	4452 UG/L	0.955	30734
745	C450	NAPHTHALENE	128	19958 UG/L	1.007	37914
858	C470	2-METHYLNAPHTHALENE	142	1196 UG/L	1.159	19280
1036	C550	ACENAPHTHENE	153	769 UG/L	1.005	6600
1062	C565	DIBENZOFURAN	168	431 UG/L	1.030	5391
1119	C590	FLUORENE	166	349 UG/L	1.083	353
275	C640	PHENANTHRENE	178	326 UG/L	1.002	3581

* SURROGATES WERE DILUTED OUT

004030

LIBRARY SEARCH
91/13/87 9:57:00 + 14:02
SAMPLE: CAV-0W02-001
COND.: —
ENHANCED (S 15B 2N 0T)

DATA: 4U12057C02A # 842 BASE M/Z: 121
CALI: 4U12057C02A # 3 RIC: 5551.

SPECTRUM DC # ----19

SAMPLE

1448

PHENOL, 3,5-DIETHYL-

C18.H14.0

M WT 158
B PK 135
RANK 1
7445
PUR 746

C10.H14.0

M WT 158
B PK 135
RANK 2
7438
PUR 642

C10.H14.0

M WT 158
B PK 135
RANK 3
7449
PUR 627

INST ID: 4530

M/Z

40

60

80

100

120

140

004030

004031

LIBRARY SEARCH
 01/13/87 9:57:00 + 15:46
 SAMPLE: CAU-DW2-001
 CONDS.: —
 ENHANCED (S 158 2N 0T)

DATA: 4U12057C02A # 946
 CALI: 4U12057C02A # 3
 BASE M/Z: 133
 RIC: 4023.

SPECTRUM DC # ---19

SAMPLE
1102

ANALYST: GU

C11.H16
 1102
 M WT 148
 B PK 133
 RANK 1
 # 7165
 PUR 668

BENZENE, PENTAMETHYL-

C10.H12.0
 1102
 M WT 148
 B PK 133
 RANK 2
 # 7154
 PUR 645

PHENOL, 2-(2-METHYL-2-PROPYL)-

C10.H12.0
 1102
 M WT 148
 B PK 133
 RANK 3
 # 7168
 PUR 643

PHENOL, P-(2-METHYLALLYL)-

INST. ID: 4530

M/Z

49 60 80 100 120 140 160

004031

004032

LIBRARY SEARCH
01/13/87 9:57:00 + 18:23
SAMPLE: CAV-0W82-801
COND.: ---
ENHANCED (S 15B 2N 8T)

DATA: 4U12057C02A #1103 BASE M/Z: 147
CALIB: 4U12057C02A # 3 RIC: 13183.

SPECTRUM DC # 19

SAMPLE

1318

ANALYST: GU

C9.H9.O.N
M WT 1318
B PK 118
RANK 1
S 7103
PUR 651

1H-INDOLE-1-CARBOXALDEHYDE, 2,3-DIHYDRO-

C7.H5.O.N3
M WT 1318
B PK 147
RANK 2
S 7682
PUR 642

PYRIDO[3,2-D]PYRIMIDIN-4-OL

C7.H5.O.N3
M WT 1318
B PK 147
RANK 3
S 7080
PUR 629

PYRIDO[3,4-D]PYRIMIDIN-4(3H)-ONE

INST ID: 4530

M/Z

40 60 80 100 120 140 160

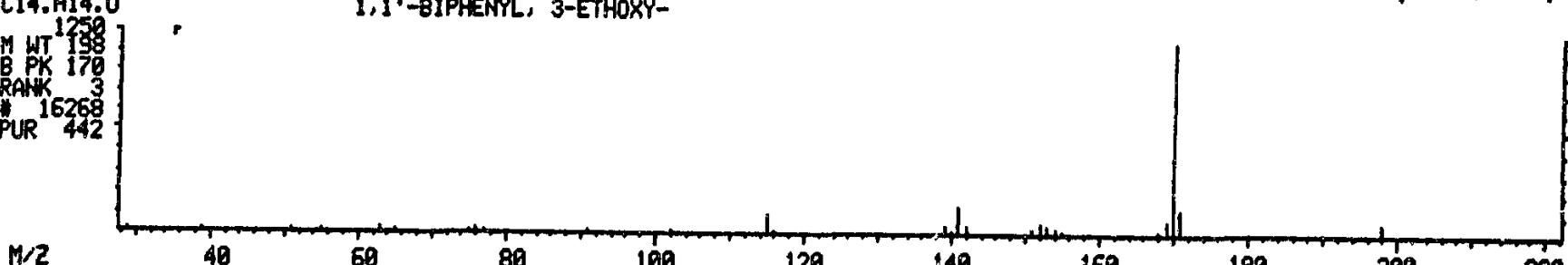
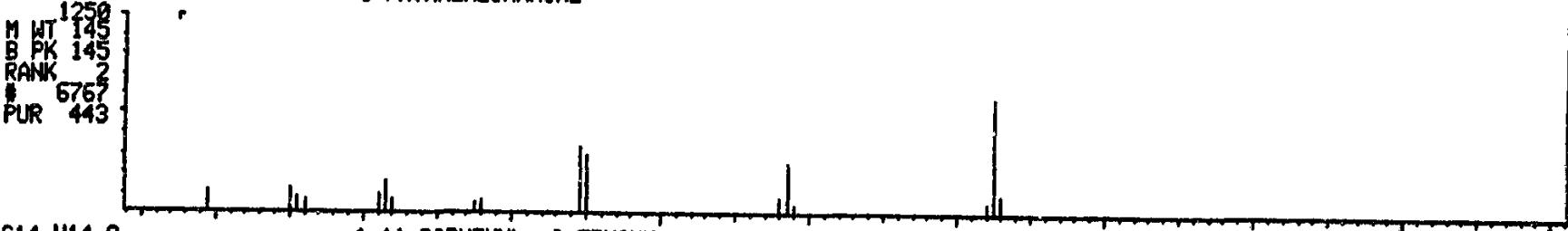
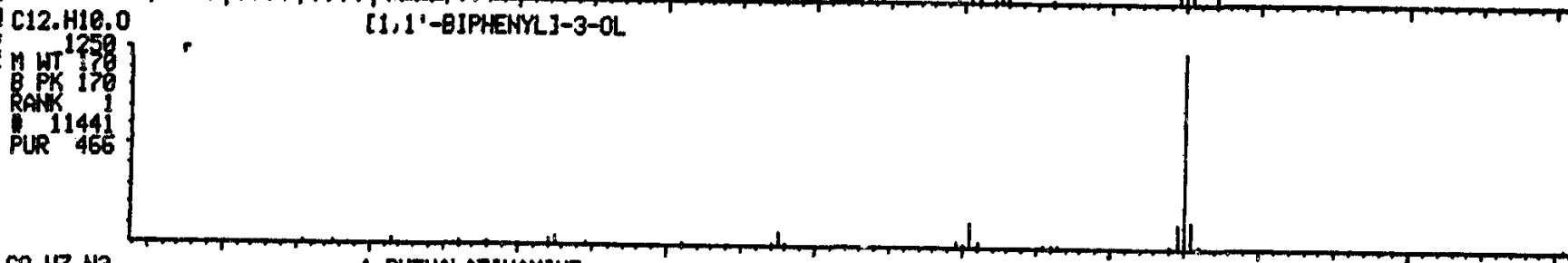
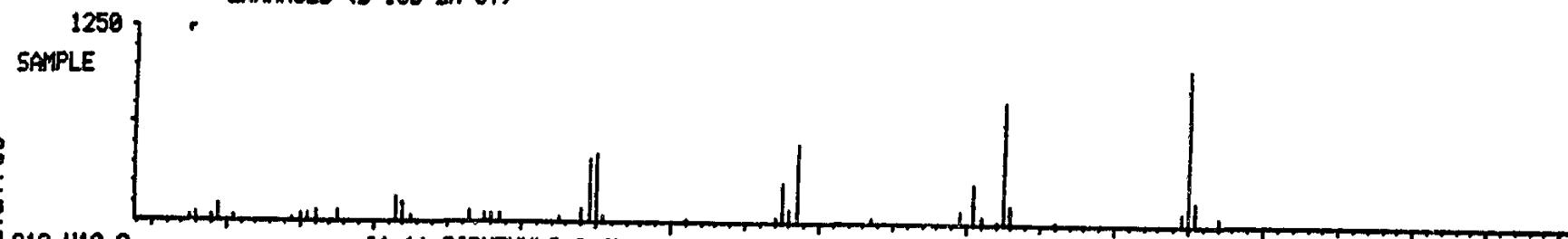
004032

004033

SPECTRUM DC # ----19

LIBRARY SEARCH
01/13/87 9:57:00 + 20:14
SAMPLE: CAV-0W02-001
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12057C02A #1214 BASE M/Z: 170
CALI: 4U12057C02A # 3 RIC: 3787.

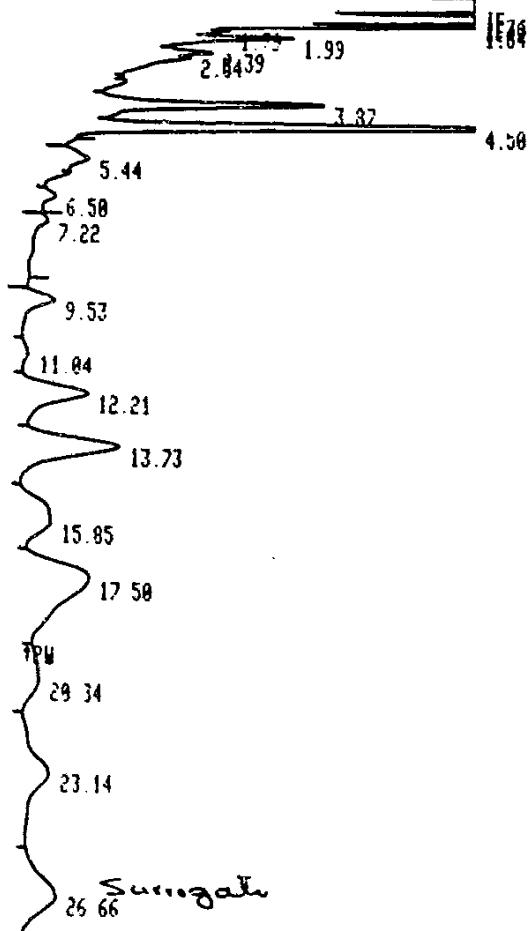


INST ID: 4530

004033

86-12-057-024

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-0W02-a
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 183 °C
VOLUME INJ. 1 μL
INSTRUMENT ID. 6000

00400

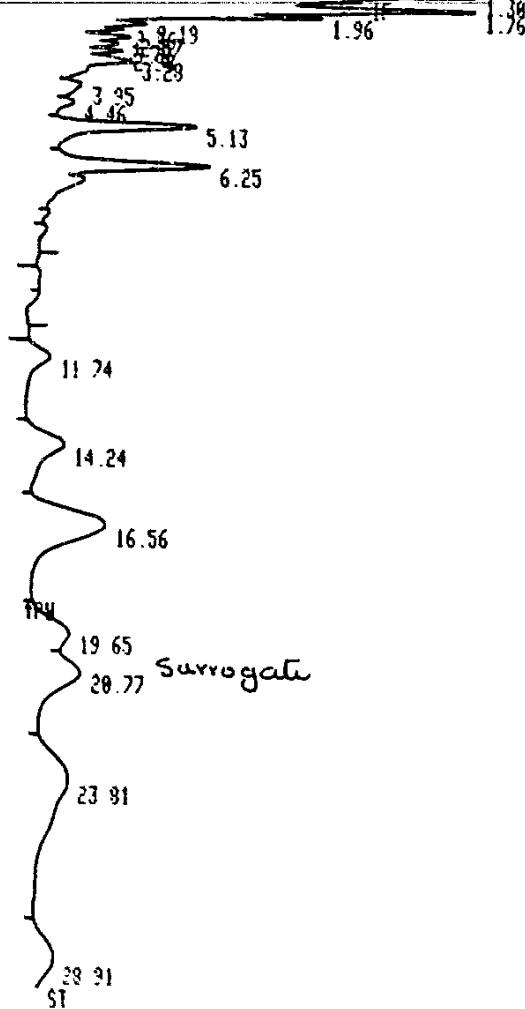
RUN # 19 JAN/86/87 22:03:28
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.36	95625	BV	0.086	14.044
1.42	141118	D VV	0.108	20.725
1.64	68691	D VP	0.083	10.088
1.79	3782	D PV	0.069	0.555
1.99	25849	D VP	0.116	3.796
2.39	13956	PV	0.145	2.050
2.54	10496	D VP	0.199	1.542
3.87	59027	PV	0.219	8.669
4.58	151073	V8	0.167	22.187
5.44	7664	BV	0.396	1.126
6.58	2883	PB	0.255	0.423
9.53	7894	BP	0.408	1.042
11.04	1595	PV	0.501	0.234
12.21	16672	VV	0.620	2.449
13.73	24977	VV	0.638	3.668
15.85	7959	VV	1.143	1.169
17.58	17589	VV	1.541	2.583
20.34	5220	VV	1.587	0.767
23.14	8089	VV	1.744	1.188
26.66	10097	V8	1.573	1.483
34.10	1451	I BH	1.320	0.213

86-12-053-02A

N/A.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-0W0Z-0C
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 203°C
VOLUME INJ. 1 μL
INSTRUMENT ID. 600

RUN # 102 JAN/12/87 14:11:14
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.30	282174	D BB	0.082	63.446
1.76	35226	VV	0.144	7.921
1.96	19581	D VP	0.102	4.403
2.19	3595	D PV	0.112	0.808
2.36	2620	D VP	0.088	0.589
2.57	4751	PV	0.142	1.068
2.78	3358	D VV	0.151	0.368
2.98	4944	VV	0.142	1.112
3.28	6706	VP	0.289	1.508
3.95	2064	PV	0.319	0.464
4.46	1827	VP	0.232	0.411
5.13	18062	PV	0.268	4.061
6.25	20450	VV	0.294	4.598
11.24	2920	BP	0.607	0.657
14.24	4869	PV	0.947	1.095
16.56	10136	VV	1.110	2.279
19.65	5453	VV	1.003	1.226
20.77	6905	VV	1.252	1.553
23.81	5256	VV	2.626	1.182
28.91	3347	I VP	1.444	0.753

Page 1
Received: 12/12/86

SPECTRO COLOR REPORT
Results by Sample

Work Order # 86-12-05

SAMPLE ID CAV-DW02-001		SAMPLE # 02 FRACTIONS: A									
		Date & Time Collected 12/12/86					Category				
AG	<10	AS	69.4	BE	<3	CD	<5	CN_SU	<100	CR	<9
	ug/L		ug/L		ug/L		ug/L		ug/L		ug/L
CU	<10	FE	4580	HG	<0.2	NI	<29	PB	<5	SB	<44
	ug/L		ug/L		ug/L		ug/L		ug/L		ug/L
SE	<5	TL	<10	ZN	29						
	ug/L		ug/L		ug/L						

CHECKLIST

SECTION II - Sample Data Packets

Sample No: CAV-0W08-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

004037

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-0W08-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205703
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *SPH*

CASE NO.: ----
 QC REPORT NO.: ----
 CONTRACT NO.: ----
 DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/19/86
 DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V03
 DILUTION FACTOR: 1:00

CAS #	COMPOUND	UG/L	8
C010	CHLOROMETHANE	10	U
C015	BROMOMETHANE	10	U
C020	VINYL CHLORIDE	10	U
C025	CHLOROETHANE	10	U
C030	METHYLENE CHLORIDE	5	U
C035	ACETONE	10	U
C040	CARBON DISULFIDE	5	U
C045	1, 1-DICHLOROETHENE	5	U
C050	1, 1-DICHLOROETHANE	5	U
C055	TRANS-1, 2-DICHLOROETHENE	5	U
C060	CHLOROFORM	5	U
C065	1, 2-DICHLOROETHANE	5	U
C110	2-BUTANONE	10	U
C115	1, 1, 1-TRICHLOROETHANE	5	U
C120	CARBON TETRACHLORIDE	5	U
C125	VINYL ACETATE	10	U
C130	BROMODICHLOROMETHANE	5	U
C140	1, 2-DICHLOROPROPANE	5	U
C145	TRANS-1, 3-DICHLOROPROPENE	5	U
C150	TRICHLOROETHENE	5	U
C155	DIBROMOCHLOROMETHANE	5	U
C160	1, 1, 2-TRICHLOROETHANE	5	U
C165	BENZENE	5	U
C170	CIS-1, 3-DICHLOROPROPENE	5	U
C175	2-CHLOROETHYL VINYLETHER	10	U
C180	BROMOFORM	5	U
C210	2-HEXANONE	10	U
C205	4-METHYL-2-PENTANONE	10	U
C220	TETRACHLOROETHENE	5	U
C225	1, 1, 2, 2-TETRACHLOROETHANE	5	U
C230	TOLUENE	5	U
C235	CHLOROBENZENE	5	U
C240	ETHYLBENZENE	5	U
C245	STYRENE	5	U
C250	TOTAL XYLENES	5	U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: CAV-DWOB-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205703
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86.

SEMOVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/18/86
 DATE ANALYZED: 01/12/87

DATAFILE: 4U12057C03
 DILUTION FACTOR: 1.0

CAS #	COMPOUND	UG/L
C315	PHENOL	10 U
C325	BIS(2-CHLOROETHYL)ETHER	10 U
C330	2-CHLOROPHENOL	10 U
C335	1,3-DICHLOROBENZENE	10 U
C340	1,4-DICHLOROBENZENE	10 U
C345	BENZYL ALCOHOL	10 U
C350	1,2-DICHLOROBENZENE	10 U
C355	2-METHYLPHENOL	10 U
C360	BIS(2-CHLOROISOPROPYL)ETHER	10 U
C365	4-METHYLPHENOL	10 U
C370	N-NITROSONDIPROPYLAMINE	10 U
C375	HEXACHLOROETHANE	10 U
C410	NITROBENZENE	10 U
C415	ISOPHORONE	10 U
C420	2-NITROPHENOL	10 U
C425	2,4-DIMETHYLPHENOL	10 U
C430	BENZOIC ACID	50 U
C435	BIS(2-CHLOROETHOXY)METHANE	10 U
C440	2,4-DICHLOROPHENOL	10 U
C445	1,2,4-TRICHLOROBENZENE	10 U
C450	NAPHTHALENE	1 U
C455	4-CHLOROANILINE	10 U
C460	HEXACHLOROBUTADIENE	10 U
C465	P-CHLORO-M-CRESOL	10 U
C470	2-METHYLNAPHTHALENE	10 U
C510	HEXACHLOROCYCLOPENTADIENE	10 U
C515	2,4,6-TRICHLOROPHENOL	10 U
C520	2,4,5-TRICHLOROPHENOL	50 U
C525	2-CHLORONAPHTHALENE	10 U
C530	2-NITROANILINE	50 U
C535	DIMETHYL PHTHALATE	10 U
C540	ACENAPHTHYLENE	10 U
C545	3-NITROANILINE	50 U
C550	ACENAPHTHENE	10 U
C555	2,4-DINITROPHENOL	50 U
C560	4-NITROPHENOL	50 U
C565	DIBENZOFURAN	10 U
C570	2,4-DINITROTOLUENE	10 U
C575	2,6-DINITROTOLUENE	10 U

004039

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-DW02-001

INMOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C03

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	10 U
C585	4-CHLOROPHENYL PHENYL ETHER	10 U
C590	FLUORENE	10 U
C595	4-NITROANILINE	50 U
C610	4, 6-DINITRO-2-METHYLPHENOL	50 U
C615	N-NITROSODIPHENYLAMINE	10 U
C625	4-BROMOPHENYL PHENYL ETHER	10 U
C630	HEXACHLOROBENZENE	10 U
C635	PENTACHLOROPHENOL	50 U
C640	PHENANTHRENE	10 U
C645	ANTHRACENE	10 U
C650	DI-N-BUTYL PHTHALATE	10 U
C655	FLUORANTHENE	10 U
C715	PYRENE	10 U
C720	BUTYL BENZYL PHTHALATE	10 U
C725	3, 3'-DICHLOROBENZIDINE	20 U
C730	BENZO(A)ANTHRACENE	10 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	4 J
C740	CHRYSENE	10 U
C760	DI-N-OCTYL PHTHALATE	10 U
C765	BENZO(B)FLUORANTHENE	10 U
C770	BENZO(K)FLUORANTHENE	10 U
C775	BENZO(A)PYRENE	10 U
C780	INDENO(1, 2, 3-CD)PYRENE	10 U
C785	DIBENZO(A, H)ANTHRACENE	10 U
C790	BENZO(QHI)PERYLENE	10 U

004040

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No : 86-12-057-03A

DC No: KEYSTONE
SMO No: CAV-DW08-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs
Conc: Low
Date Extracted: 12/18/86
Date Analyzed: 1/6/87
Dilution Factor: 1
Percent Moisture(decant):
GFC Cleanup Yes ✓
Sep. Funnel Ext. Yes ✓
Cont. Liquid Ext. Yes ✓

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.05 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYPHENYL CHLOR	0.10 U
53494-70-5	ENDRIN KETONE	0.50 U
57-74-9	CHLORDANE	0.10 U
8001-35-2	TOXAPHENE	0.50 U
12674-11-2	AROCLOTR-1016	1.0 U
11104-28-2	AROCLOTR-1221	0.50 U
11141-16-5	AROCLOTR-1232	0.50 U
53469-21-9	AROCLOTR-1242	0.50 U
12672-29-6	AROCLOTR-1248	0.50 U
11097-69-1	AROCLOTR-1254	0.50 U
11096-82-5	AROCLOTR-1260	1.0 U
		1.0 U

V_s = 500 = Volume of water extracted (ml)
V_t = 5000 = Volume of total extract (ul)
V_i = 1 = Volume of extract injected(ul)
1st Column Inst.: 6000
2nd Column Inst.: 6000 NA

U=UNDETECTED AT THE LISTED DETECTION LIMIT.
J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.
B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: CAV-0W08-00

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: NOT

DATAFILE: 4U12057V03

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
-------	-------------------------	-------	--------	--------

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

004042

INST ID: 4530

SPECTRIX DG # ---- 8

SAMPLE: CAV-QW08

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12057C03

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
110-12-3	2-HEXANONE, 5-METHYL-	* 296	926	15 43
25144-05-2	CYCLOPENTANOL, 2-METHYL-, CIS-	337	893	10 40
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	761		10 Q
4381-40-1	PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHY	1129	915	6 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1210		10 J
112-24-9	ATRAZINE (ACN)	1249	785	9 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

004044

RIC
12/19/86 10:40:00
SAMPLE: CAV-0W08-001
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057V03 #1
CALI: 4U12057V03 #3

SCANS 25 TO 600

146176.

SPECTRIX DC # -----14

100.0

RIC

INST ID: 4000

TS1

173

SM1

231

TS2

386

TS2

459

TS3

482

TS3

555

100

5:00

200

10:00

300

15:00

400

20:00

500

25:00

600 SCAN
20:00 TIME

004044

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID.

CAV-0W08-001

FILENAME

4U12057V03

CLIENT

KEYSTONE

INST

4000

DATE INJECTED

12/19/86 10:40:00

ANALYST

NOT

CALIB STD DATE

12-14-86

VERIFIED BY

NOT

CORR. FACTOR

1.00

SCAN#	VOA COMPOUNDS		M/E	AMOUNT	RRT	AREA
173 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000	15952
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000	73509
482 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000	63659
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	92 %	1.335	21750
459 CS05	D8-TOLUENE	**SU2**	98	99 %	0.952	64444
555 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	94 %	1.151	44050
110 C030	METHYLENE CHLORIDE *		84	3 UG/L	0.636	1760

* BLANK CORRECTED ON DATA SHEET

004046

RIC
01/12/87 16:58:00
SAMPLE: CAV-DW88-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057C03 #1
CALI: 4U12057C03 #3
SCANS 200 TO 1350

114568.

SPECTRUM DC # ----14

INST ID: 4530

100.0

RIC

SU-1

SU-2

SU-3

SU-4

SU-5

TS-1

TS-2

TS-3

TS-4

TS-5

200

31

400

6:40

600

10:00

700

7:20

800

7:20

1000

16:40

1200

28:00

SCAN

TIME

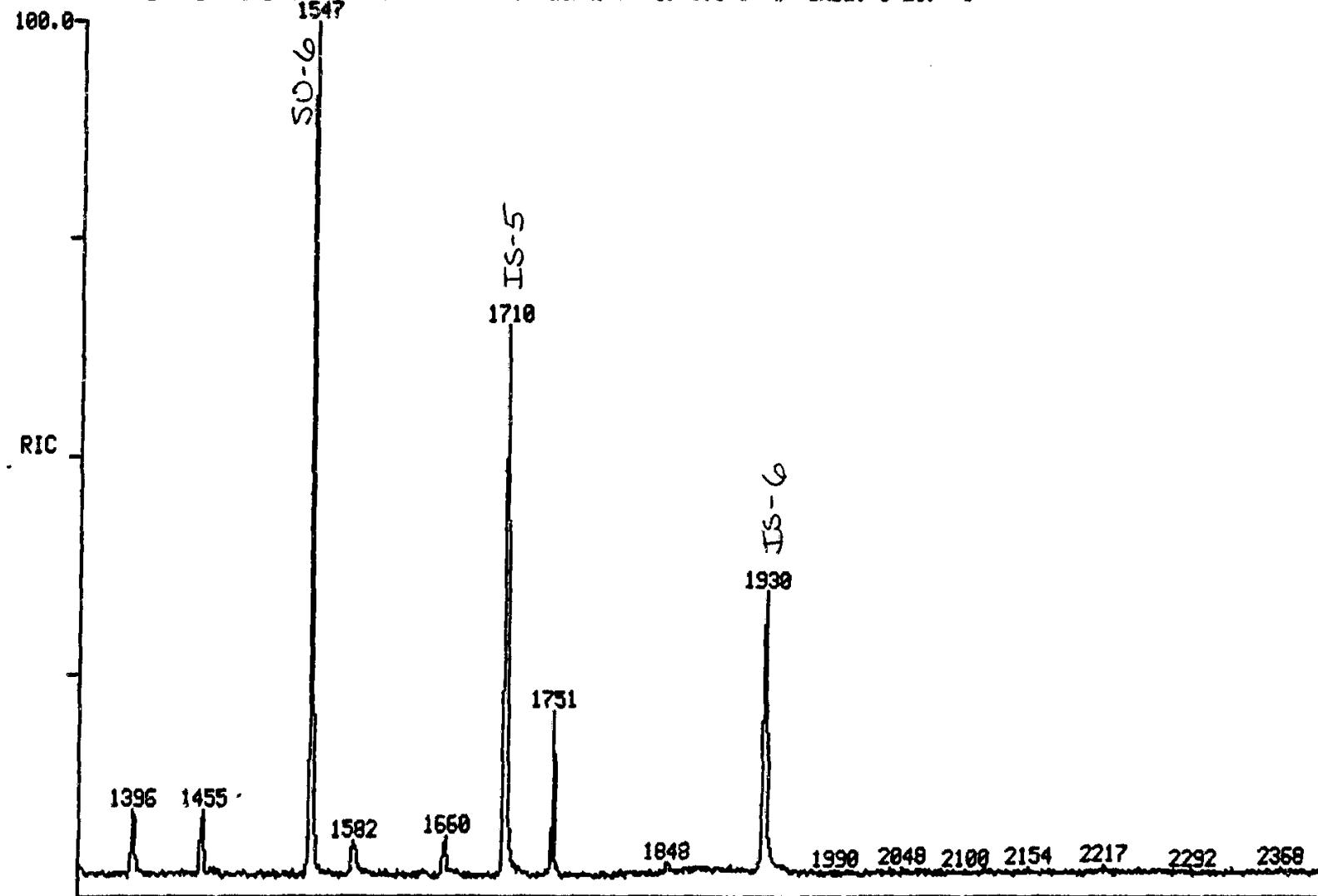
004046

004047

RIC
01/12/87 16:58:00
SAMPLE: CAV-OM08-001
COND.S.: ---
RANGE: G 1.2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057C03 #1
CALIB: 4U12057C03 #3
SCANS 1350 TO 2400

SPECTRIX DC # 14



INST ID: 4530

1400 23120 1600 26140 1800 30100 2000 33120 2200 36140 2400 36140 SCAN
--:00 TIME

004047

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	CAV-0W08-001		
FILENAME	4U12057C03	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/12/87 16:58:00	VERIFIED BY	<i>JPM</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	1.00

SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
325	CI30	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	40 UG/L	1.000	2011:
736	CI40	NAPHTHALENE-D8 ** IS2 **	136	40 UG/L	1.000	7479:
1030	CI50	ACENAPHTHENE-D10 ** IS3 **	164	40 UG/L	1.000	3593:
1270	CI60	PHENANTHRENE-D10 ** IS4 **	188	40 UG/L	1.000	4322:
1710	CI70	CHRYSENE-D12 ** IS5 **	240	40 UG/L	1.000	2619:
1930	CI75	PERYLENE-D12 ** IS6 **	264	40 UG/L	1.000	1713:
329	CS50	2-FLUOROPHENOL ** SU1 **	112	51 %	0.627	3609:
493	CS45	PHENOL-D5 ** SU2 **	99	56 %	0.939	5885:
1161	CS55	2,4,6-TRIBROMOPHENOL ** SU5 **	330	55 %	1.127	5821:
622	CS20	NITROBENZENE-D5 ** SU3 **	82	102 %	0.845	4463:
927	CS25	2-FLUOROBIPHENYL ** SU4 **	172	98 %	0.900	9150:
1547	CS30	TERPHENYL-D14 ** SU6 **	244	85 %	0.905	3031:
739	C450	NAPHTHALENE	128	1 UG/L	1.004	192
1751	C735	BIS(2-ETHYLHEXYL)PHTHALATE	149	4 UG/L	1.024	282

004049

LIBRARY SEARCH
 01/12/87 16:58:00 + 12:41
 SAMPLE: CAV-DW08-001
 CONDS.: --
 ENHANCED (S 15B 2N 0T)

DATA: 4U12057C03 # 761 BASE M/Z: 81
 CALI: 4U12057C03 # 3 RIC: 13167.

SPECTRIX DC # ----19

ANALYST: GU

C8.H16.0

SAMPLE

M WT 1000
 B PK 128
 RANK 1
 # 4092
 PUR 724

FURAN, 2,5-DIETHYLtetrahydro-

C12.H22.02

M WT 1000
 B PK 198
 RANK 2
 # 16210
 PUR 659

2,2'-BIOXEPANE

C10.H18.0

M WT 1000
 B PK 154
 RANK 3
 # 8276
 PUR 629

CYCLOHEXENE, 3-(2-METHYLPROPOXY)-

INST ID: 4530

M/Z

40 60 80 100 120 140 160 180 200 220

004049

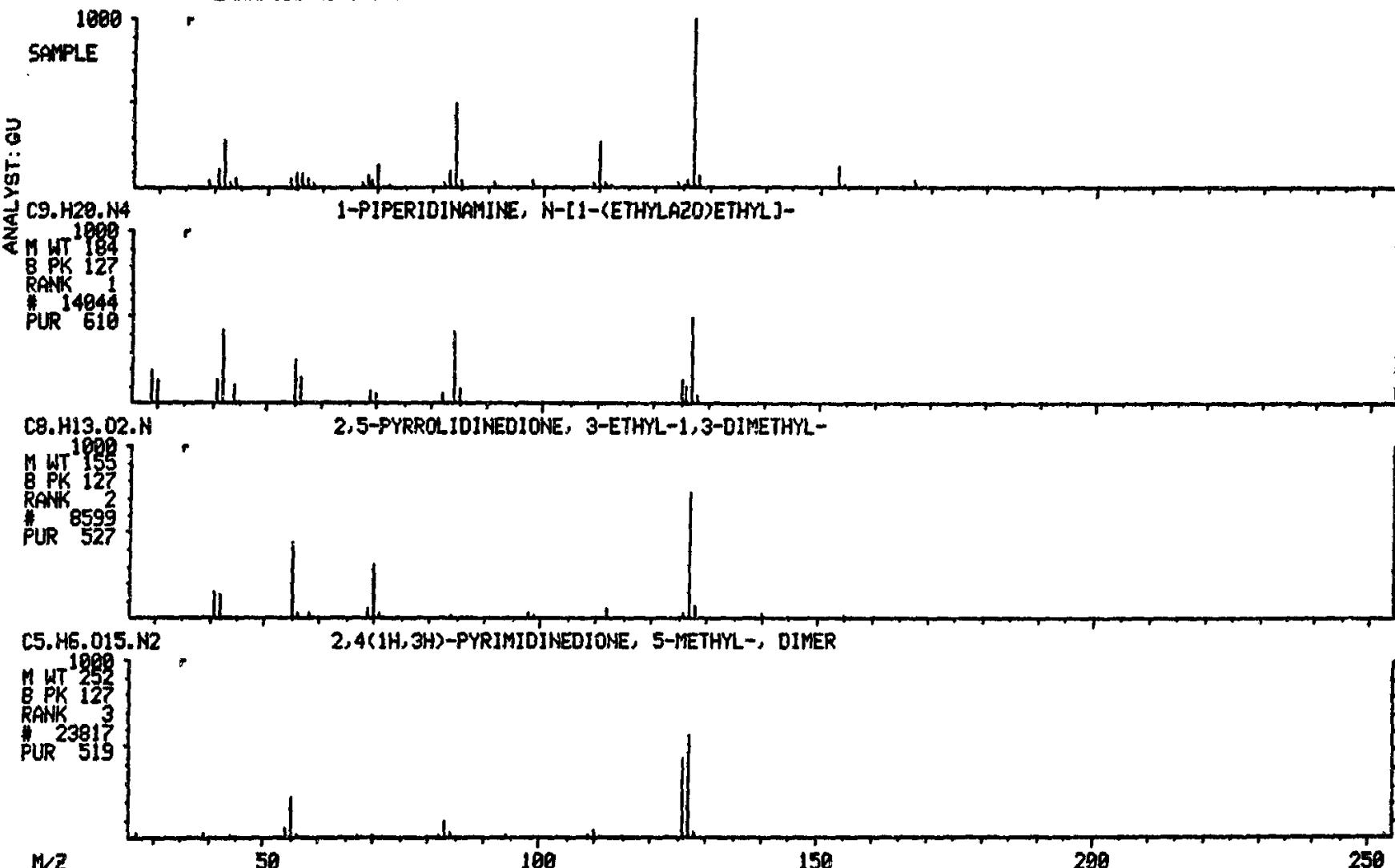
SPECTRUM DC # -----19

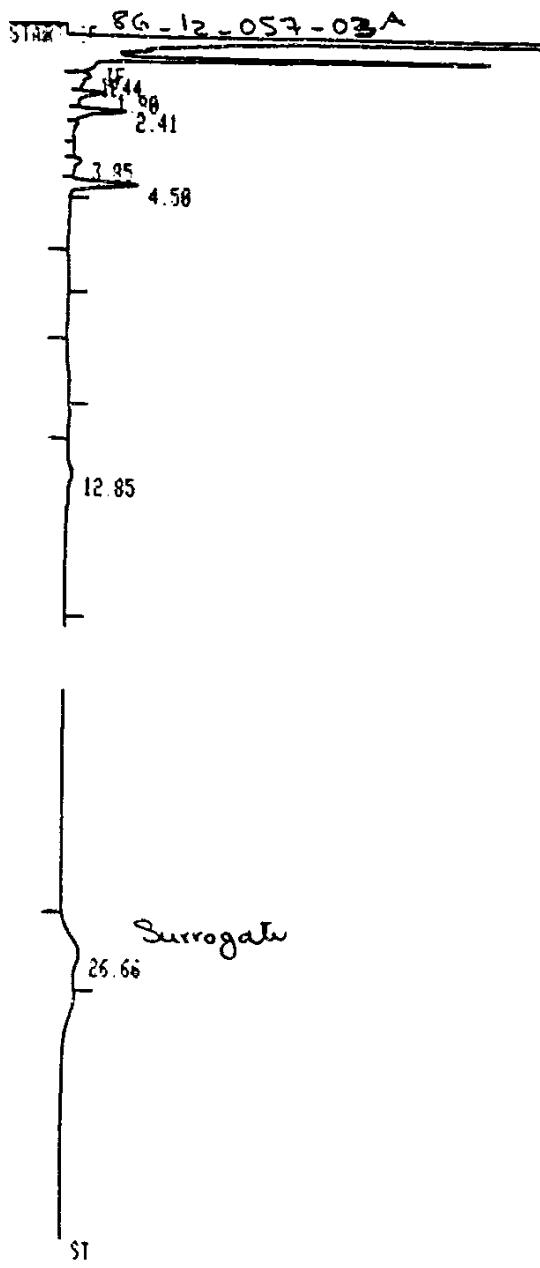
INST ID: 4330

004050

LIBRARY SEARCH
01/12/87 16:58:00 + 20:10
SAMPLE: CAV-0408-001
COND.: ----
ENHANCED (S 15B 2N 0T)

DATA: 4U12057C03 #1210
CALI: 4U12057C03 # 3
BASE M/Z: 127
RIC: 5215.





DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-0W08-00
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 μL
INSTRUMENT ID. 6000

004
00

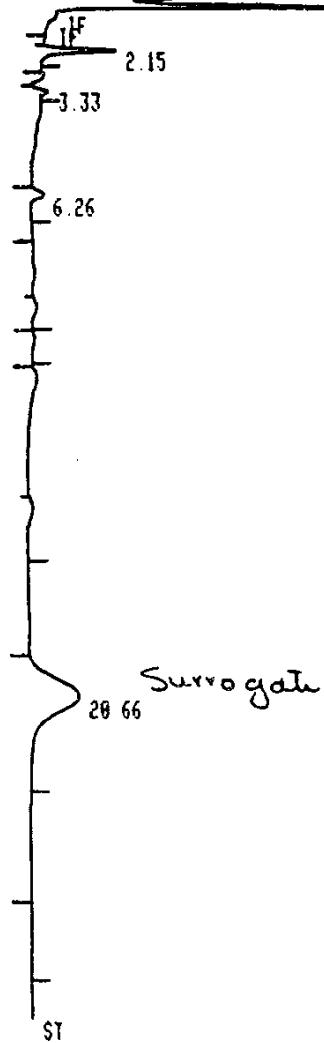
RUN # 20
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.44	1866	PP	0.201	4.176
1.98	6791	PP	0.120	15.198
2.41	12504	PP	0.137	27.984
3.85	2121	PV	0.233	4.747
4.50	17177	VB	0.123	38.442
12.85	1511	BB	0.764	3.382
26.66	2713	PB	0.898	6.072

STRTLF 86-12-057-03A

NA



DC# _____
SPECTRIX HOUSTON

CASE NO. Key Stone

SAMPLE NO. CAV-0W08-O
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 203°C
VOLUME INJ. 1 uL
INSTRUMENT ID. 69C

0
0

RUN # 103 JAN/12/87 14:42:31
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%			
RT	HEIGHT	TYPE	AR/HT
2.15	9192	PB	0.142
3.33	1526	PB	0.178
6.26	1660	BB	0.228
20.66	6383	PB	1.172

TOTAL HGT= 13761
MUL FACTOR= 1.0000E+00

Page 1

Received: 12/12/86

SPECTRUM 00

REPORT
Results by Sample

Work Order # 86-12-05

SAMPLE ID CAV-DW08-001		SAMPLE # 03 FRACTIONS: A		Category	
Date & Time Collected 12/12/86					
AG	<10	AS	22.2	BE	<3
	ug/L		ug/L		ug/L
CD	9	CN	SU	<100	CR
	ug/L			ug/L	ug/L
CU	13	FE	5340	HG	<0.2
	ug/L		ug/L		ug/L
NI	<29	PB	27.8	SB	<44
	ug/L		ug/L		ug/L
SE	<5	TL	<10	ZN	48
	ug/L		ug/L		ug/L

CHECKLIST

SECTION II - Sample Data Packets

- Sample No: CAV-0W09-001
- Organic Analysis Data Sheets (OADS)
 - VOA
 - SV
 - Pest.
 - Tentatively Identified Compounds
 - VOA
 - SV
 - Raw Data
 - VOA chromatogram
 - VOA quantitation report (surrogates)
 - SV chromatogram
 - SV quantitation report (surrogates)
 - Pest. chromatogram and integrator print out
 - Inorganic Analysis Data Sheets (IADS)

004054

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-0W09-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205704
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *YDH*

CASE NO.: -----
QC REPORT NO.: -----
CONTRACT NO.: -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V04
DILUTION FACTOR: 1.00

CAS #	COMPOUND	UG/L.	5
C010	CHLOROMETHANE	10 U	
C015	BROMOMETHANE	10 U	
C020	VINYL CHLORIDE	10 U	
C025	CHLOROETHANE	10 U	
C030	METHYLENE CHLORIDE	5 U	
C035	ACETONE	10 U	
C040	CARBON DISULFIDE	5 U	
C045	1, 1-DICHLOROETHENE	5 U	
C050	1, 1-DICHLOROETHANE	5 U	
C055	TRANS-1, 2-DICHLOROETHENE	5 U	
C060	CHLOROFORM	5 U	
C065	1, 2-DICHLOROETHANE	5 U	
C110	2-BUTANONE	10 U	
C115	1, 1, 1-TRICHLOROETHANE	5 U	
C120	CARBON TETRACHLORIDE	5 U	
C125	VINYL ACETATE	10 U	
C130	BROMODICHLOROMETHANE	5 U	
C140	1, 2-DICHLOROPROPANE	5 U	
C145	TRANS-1, 3-DICHLOROPROPENE	5 U	
C150	TRICHLOROETHENE	5 U	
C155	DIEBROMOCHLOROMETHANE	5 U	
C160	1, 1, 2-TRICHLOROETHANE	5 U	
C165	BENZENE	5 U	
C170	CIS-1, 3-DICHLOROPROPENE	5 U	
C175	2-CHLOROETHYL VINYLETHER	10 U	
C180	BROMOFORM	5 U	
C210	2-HEXANONE	10 U	
C205	4-METHYL-2-PENTANONE	10 U	
C220	TETRACHLOROETHENE	5 U	
C225	1, 1, 2-TETRACHLOROETHANE	5 U	
C230	TOLUENE	5 U	
C235	CHLOROBENZENE	5 U	
C240	ETHYLBENZENE	5 U	
C245	STYRENE	5 U	
C250	TOTAL XYLENES	5 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: CAV-DW09-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205704
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *Y.P.H.*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86.

SEMIVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: ... 12/12/86.
 DATE ANALYZED: 01/12/87

DATAFILE: 4U12057C04

DILUTION FACTOR: ... 1:0

CAS #	COMPOUND	UG/L	004056
C315	PHENOL	10 U	
C325	BIS(2-CHLOROETHYL)ETHER	10 U	
C330	2-CHLOROPHENOL	10 U	
C335	1, 3-DICHLOROBENZENE	10 U	
C340	1, 4-DICHLOROBENZENE	10 U	
C345	BENZYL ALCOHOL	10 U	
C350	1, 2-DICHLOROBENZENE	10 U	
C355	2-METHYLPHENOL	10 U	
C360	BIS(2-CHLOROISOPROPYL)ETHER	10 U	
C365	4-METHYLPHENOL	10 U	
C370	N-NITROSODIISOPROPYLAMINE	10 U	
C375	HEXAChLORoETHANE	10 U	
C410	NITROBENZENE	10 U	
C415	ISOPHORONE	10 U	
C420	2-NITROPHENOL	10 U	
C425	2, 4-DIMETHYLPHENOL	10 U	
C430	BENZOIC ACID	50 U	
C435	BIS(2-CHLOROETHOXY)METHANE	10 U	
C440	2, 4-DICHLOROPHENOL	10 U	
C445	1, 2, 4-TRICHLOROBENZENE	10 U	
C450	NAPHTHALENE	10 U	
C455	4-CHLOROANILINE	10 U	
C460	HEXAChLOROBUTADIENE	10 U	
C465	P-CHLORO-M-CRESOL	10 U	
C470	2-METHYLNAPHTHALENE	10 U	
C510	HEXAChLOROCYCLOPENTADIENE	10 U	
C515	2, 4, 6-TRICHLOROPHENOL	10 U	
C520	2, 4, 5-TRICHLOROPHENOL	50 U	
C525	2-CHLORONAPHTHALENE	10 U	
C530	2-NITROANILINE	50 U	
C535	DIMETHYL PHTHALATE	10 U	
C540	ACENAPHTHYLENE	10 U	
C545	3-NITROANILINE	50 U	
C550	ACENAPHTHENE	10 U	
C555	2, 4-DINITROPHENOL	50 U	
C560	4-NITROPHENOL	50 U	
C565	DIBENZOFURAN	10 U	
C570	2, 4-DINITROTOLUENE	10 U	
C575	2, 6-DINITROTOLUENE	10 U	

SPECTRIX DC # --- 8

SAMPLE NUMBER: CAV-QW09-001

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C04

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	10 U
C585	4-CHLOROPHENYL PHENYL ETHER	10 U
C590	FLUORENE	10 U
C595	4-NITROANILINE	50 U
C610	4, 6-DINITRO-2-METHYLPHENOL	50 U
C615	N-NITROSODIPHENYLAMINE	10 U
C625	4-BROMOPHENYL PHENYL ETHER	10 U
C630	HEXACHLOROBENZENE	10 U
C635	PENTACHLOROPHENOL	50 U
C640	PHENANTHRENE	10 U
C645	ANTHRACENE	10 U
C650	DI-N-BUTYL PHTHALATE	10 U
C655	FLUORANTHENE	10 U
C715	PYRENE	10 U
C720	BUTYL BENZYL PHTHALATE	10 U
C725	3, 3'-DICHLOROBENZIDINE	20 U
C730	BENZO(A)ANTHRACENE	10 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	2 J
C740	CHRYSENE	10 U
C760	DI-N-OCTYL PHTHALATE	10 U
C765	BENZO(B)FLUORANTHENE	10 U
C770	BENZO(K)FLUORANTHENE	10 U
C775	BENZO(A)PYRENE	10 J
C780	INDENO(1, 2, 3-CD)PYRENE	10 U
C785	DIBENZO(A, H)ANTHRACENE	10 U
C790	BENZO(GH)PERYLENE	10 U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

004057

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-04A

DC No: KEYSTONE
SMO No: CAV-DW09-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low

GPC Cleanup Yes No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

Date Extracted: 12/18/86
Date Analyzed: 1/6/87

Dilution Factor: 1

Percent Moisture (decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOL-1016	0.50 U
11104-28-2	AROCLOL-1221	0.50 U
11141-16-5	AROCLOL-1232	0.50 U
53469-21-9	AROCLOL-1242	0.50 U
12672-29-6	AROCLOL-1248	0.50 U
11097-69-1	AROCLOL-1254	1.0 U
11096-82-5	AROCLOL-1260	1.0 U

004050

V_s = 500 = Volume of water extracted (ml)
V_t = 5000 = Volume of total extract (ul)
V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000
2nd Column Inst. : 6000 NA

U=UNDETECTED AT THE LISTED DETECTION LIMIT.
J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.
B=COMPOUND ALSO FOUND IN BLANK.

Form 1

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: CAV-OW09-00

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: NOT

DATAFILE: 4U12057V04

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
-------	-------------------------	-------	--------	--------

004059
004059
004059
004059

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4930

SPECTRIX DG # ---- 8

SAMPLE: CAV-DW09 - 001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12057C04

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
110-12-3	2-HEXANONE, 5-METHYL-	*325	906	94 J
144-19-4	1,3-PENTANEDIOL, 2,2,4-TRIMETHYL-	706	917	48 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	764		28 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	834		7 J
-4381-40-1	PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHY	1129	902	26 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1175		14 J
207-65-9	BENZENESULFONAMIDE, N-BUTYL-4-METHYL-	1221	777	11 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1583		24 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID.	CAV-0W09-001	INST	400C
FILENAME	4U12057V04	ANALYST	NOT
CLIENT	KEYSTONE	VERIFIED BY	<i>NOT</i>
DATE INJECTED	12/19/86 11:21:00	CORR. FACTOR	1
CALIB STD DATE	12-14-86		

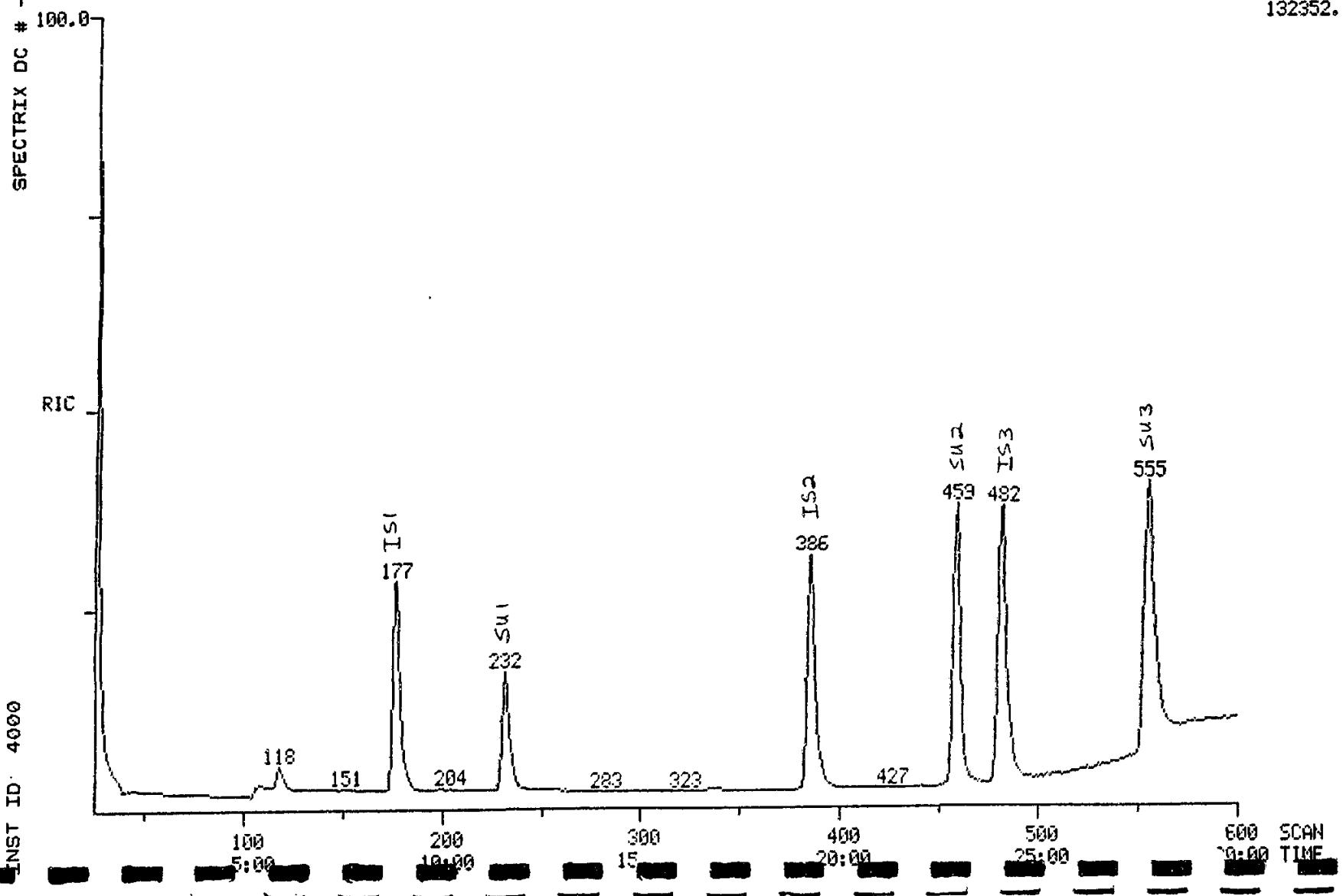
SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT
176 CJ01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L 1.000
385 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L ±.000
481 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L ±.000
232 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	92 % 1.318
459 CS05	D8-TOLUENE	**SU2**	98	99 % 0.954
555 CS10	F-BROMOFLUOROBENZENE	**SU3**	95	96 % 1.154
118 C030	METHYLENE CHLORIDE*		84	3 UG/L 0.676

* BLANK CORRECTED ON DATA SHEET

RIC
12/13/86 11:21:00
SAMPLE: CAV-0W03-001
COND.: ----
RANGE: G 1, 600 L

Z 9 0 4 0 6 2
DATA: 4U12057V04 #1 SCANS 25 TO 600
CALI: 4U12057V04 #3
0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

132352.



004062

INST ID: 4530

SPECTRIX DC # ---14



RIC
01/12/87 18:18:00
SAMPLE: CAV-ON09-001
COND.S.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

E 907.00
DATA: 4U12057C04 #1
CALI: 4U12057C04 #3
SCANS 200 TO 1350

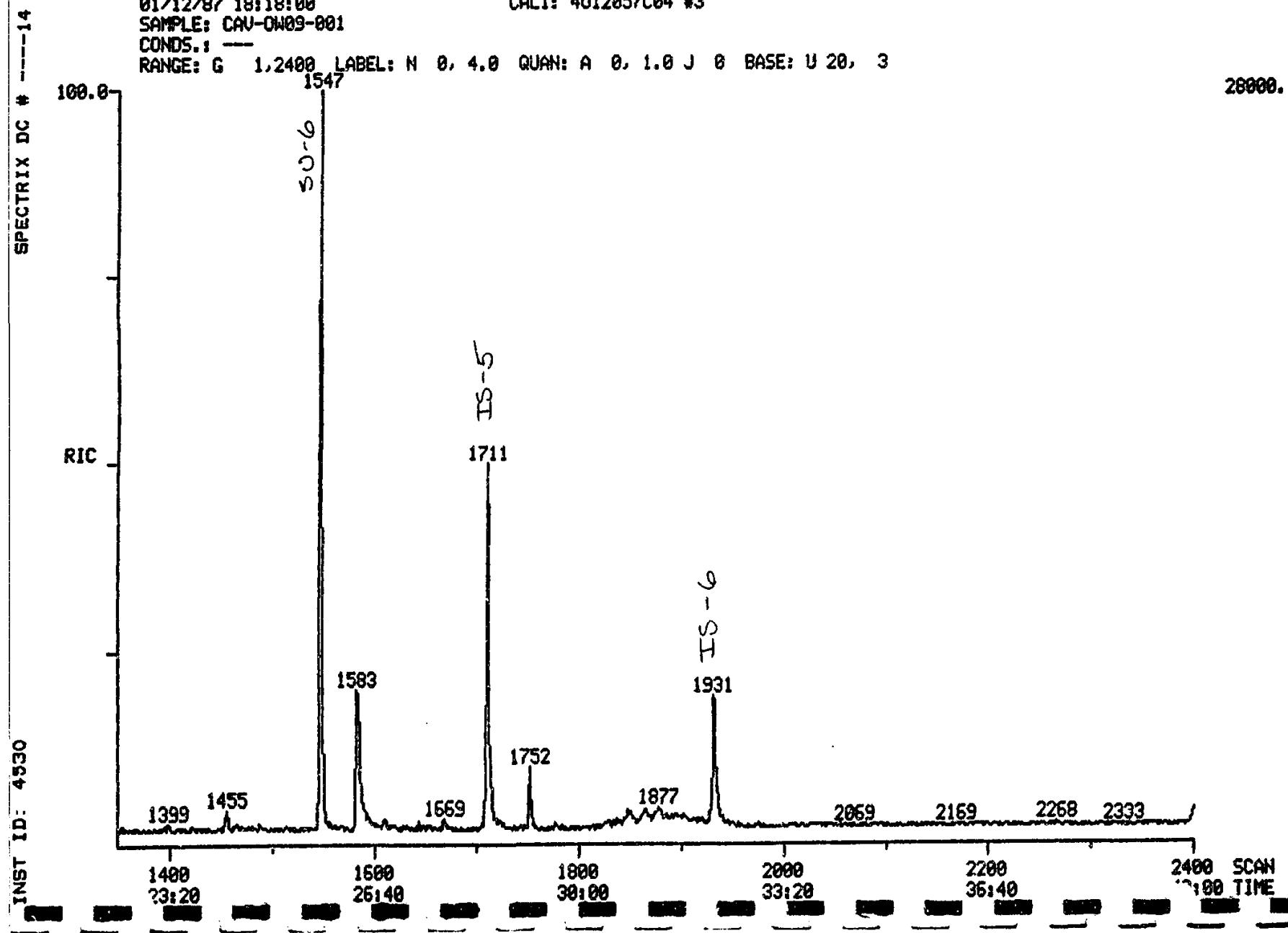
98384.

004063

SCAN
TIME

RIC
01/12/87 18:18:00
SAMPLE: CAV-0W09-001
COND.S.: —
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

79 DATA: 4U12057C04 #1
CALI: 4U12057C04 #3
SCANS 1350 TO 2400



SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	CAV-DW09-001		
FILENAME	4U12057C04	INST ID.	4930
CLIENT	KEYSTONE	ANALYST	QU
DATE INJECTED	01/12/87 18:18:00	VERIFIED BY	<i>YPH</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	1.00

SCAN#	SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
536	CI30 1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	40 UG/L	1.000	17754
739	CI40 NAPHTHALENE-D8 ** IS2 **	136	40 UG/L	1.000	68101
1030	CI50 ACENAPHTHENE-D10 ** IS3 **	164	40 UG/L	1.000	32271
1270	CI60 PHENANTHRENE-D10 ** IS4 **	188	40 UG/L	1.000	40516
1711	CI70 CHRYSENE-D12 ** IS5 **	240	40 UG/L	1.000	149914
1931	CI75 PERYLENE-D12 ** IS6 **	264	40 UG/L	1.000	9934
353	CS50 2-FLUOROPHENOL ** SU1 **	112	72 %	0.659	45004
506	CS45 PHENOL-D5 ** SU2 **	99	79 %	0.944	72431
1161	CS55 2, 4, 6-TRIBROMOPHENOL ** SU5 **	330	77 %	1.127	7514
629	CS20 NITROBENZENE-D5 ** SU3 **	82	103 %	0.851	Q0804
927	CS25 2-FLUOROBIPHENYL ** SU4 **	172	96 %	0.900	Q9971
1547	CS30 TERPHENYL-D14 ** SU6 **	244	113 %	0.904	30511
1752	C735 BIS(2-ETHYLHEXYL)PHTHALATE	149	2 UG/L	1.024	1241

INST ID: 4530

SPECTRUM DC # -----19

ANALYST: GU

SAMPLE
1000

LIBRARY SEARCH
01/12/87 18:18:00 + 12:44
SAMPLE: CAV-0W09-901
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12057C04 # 764 BASE M/Z: 81
CALI: 4U12057C04 # 3 RIC: 35583.

C8.H16.0
M WT 1099
B PK 128
RANK 1
4092
PUR 729

FURAN, 2,5-DIETHYLtetrahydro-

C12.H22.02
M WT 198
B PK 99
RANK 2
16216
PUR 657

2,2'-BIOXEPANE

C10.H18.0
M WT 1000
B PK 154
RANK 3
8270
PUR 626

CYCLOHEXENE, 3-(2-METHYLPROPOXY)-

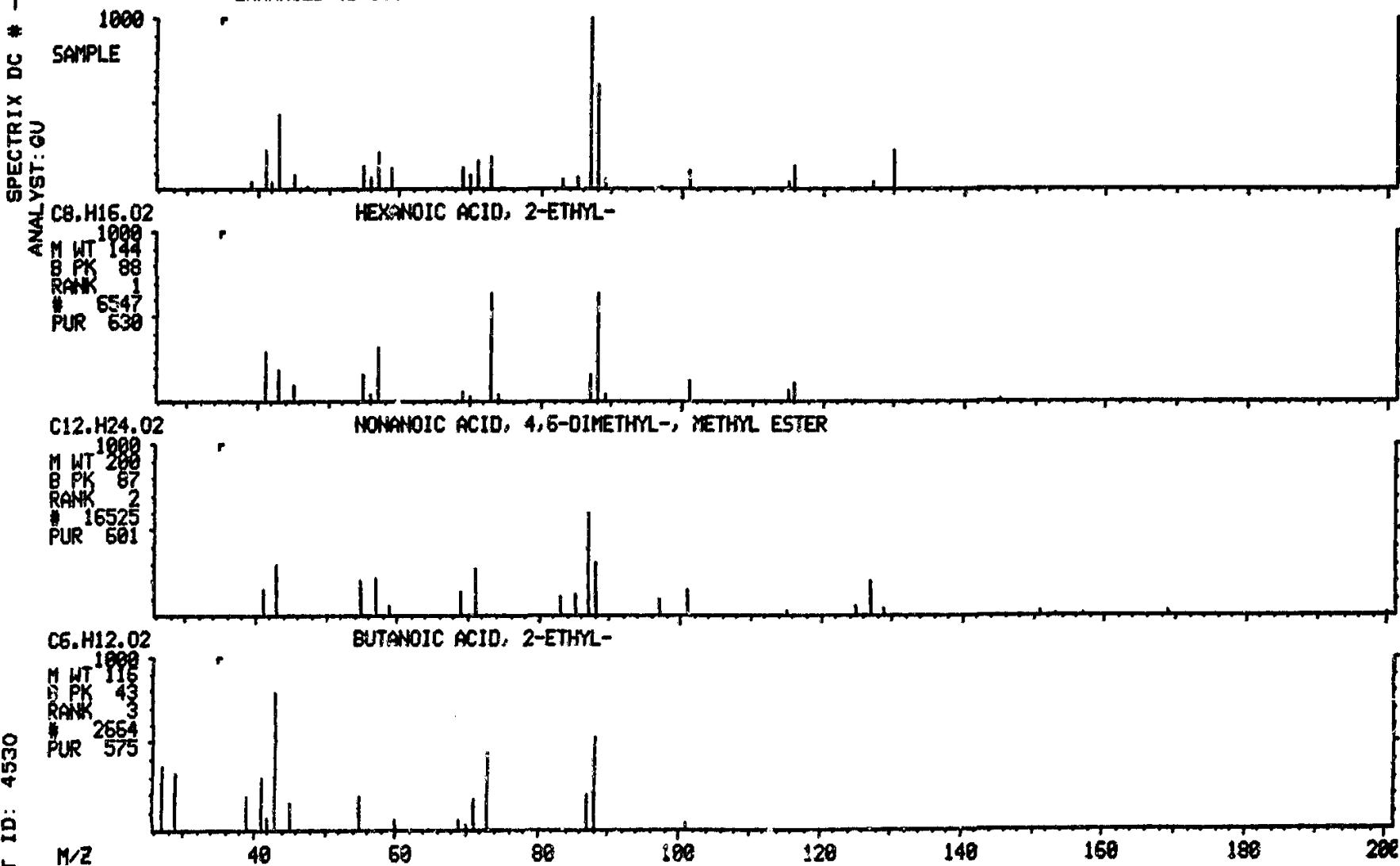
M/Z 40 60 80 100 120 140 160 180 200 220

SPECTRUM DC # ---19

INST ID: 4530

LIBRARY SEARCH
01/12/87 18:18:00 + 13:54
SAMPLE: CAN-OH09-001
CONDNS.: ---
ENHANCED (S 158 2N 0T)

L 9 0 7 0 0
DATA: 4U12857C04 # 834
CALI: 4U12857C04 # 3
BASE M/Z: 87
RIC: 2887.

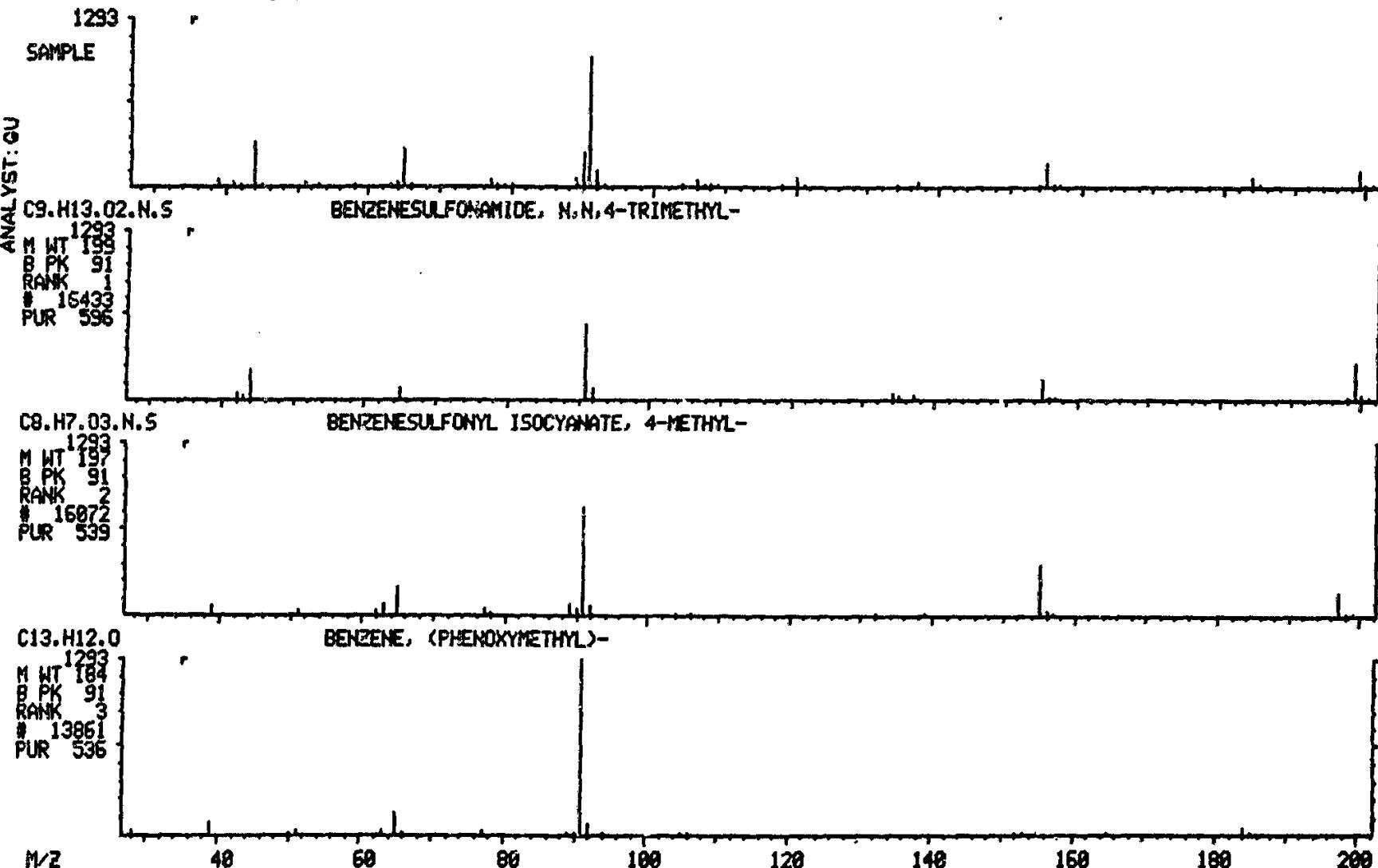


INST ID: 4530

SPECTRUM DC # -----19

LIBRARY SEARCH
01/12/87 19:18:00 + 19:35
SAMPLE: CAU-DW03-001
COND.: ----
ENHANCED (S 158 2N 0T)

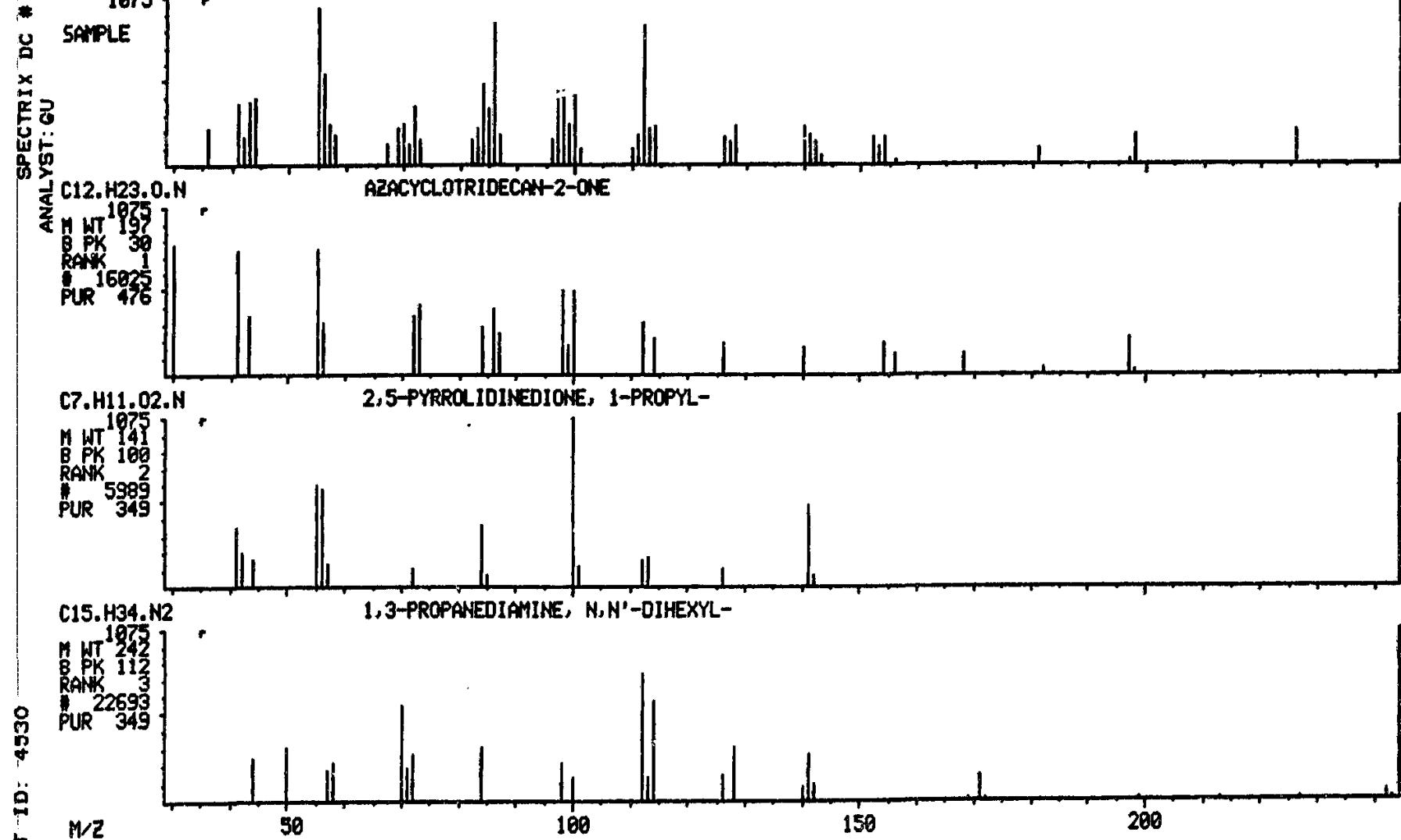
DATA: 4U12057C04 #1175
CALI: 4U12057C04 # 3
BASE M/Z: 91
RICI: 9151.



SPECTRUM DC # ---19

LIBRARY SEARCH 6 9 0 4 0 0
01/12/87 18:18:00 + 26:23
SAMPLE: CAV-0W09-001
COND.:
ENHANCED (S 15B 2N 8T)

DATA: 4U12057C04 #1583 BASE M/Z: 55
CALI: 4U12057C04 # 3 RIC: 4223.



1986-12-053-04A

NA.

IE43

3.22

DC#
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-0W09-00
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C

1.8m x2mm ID x6mm OD
3% SP-2100

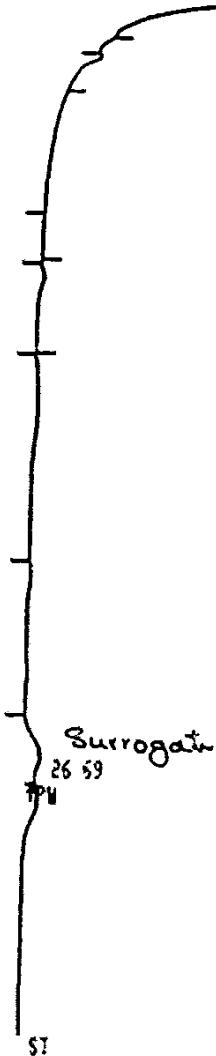
100/120 SUPELCORPRT

COLUMN TEMP. 187 °C

VOLUME INJ. 1 μl

INSTRUMENT ID. G600

0
0



RUN # 21
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

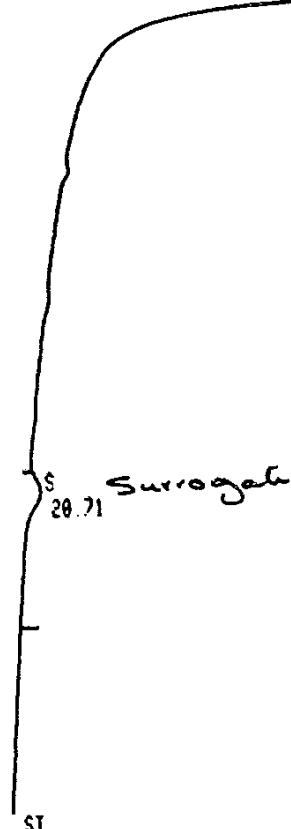
RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.43	349835	D PY	0.091	9.394
3.22	3370120	VB	2.033	90.493
26.59	4204	BY	1.161	0.113

TOTAL HC4T= 3724200
MUL FACTOR= 1.0000E+00

86-12-057-04A

NA.

F₃₄
F₃₆
F₃₇
F₃₉
4.19
6.27



DC# _____
SPECTRIX HOUSTON

CASE NO. Key Stone

SAMPLE NO. CAV.OW09-O
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm C
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 303°C
VOLUME INJ. 1µl
INSTRUMENT ID. 5600

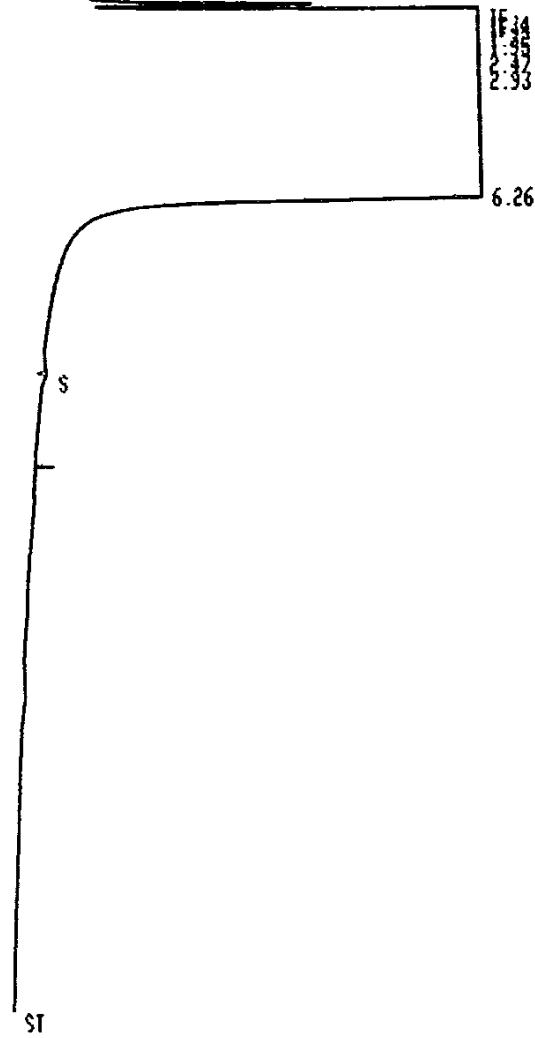
0
0

RUN # 104 JAN/12/87 15:13:47
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%	RT	HEIGHT	TYPE	AR/HT	HEIGHT%
	1.34	1016675	SPB	0.169	8.882
	1.96	1509975	TBV	0.358	13.191
	2.18	1354546	TVV	0.162	11.833
	2.47	1928583	TVV	0.459	16.848
	2.91	1358051	DTVV	0.866	11.864
	4.19	781580	TVV	0.281	6.828
	6.27	3496216	TVB	0.962	30.542
	20.71	1533	BB	0.948	0.013

TOTAL HGT= 1.1447E+02
MUL FACTOR= 1.0000E+00

STARFILE 86-12-057-04A 1:2 NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Key Stone

SAMPLE NO CAV-0W09-O
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x 6mm I
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 50.3°C
VOLUME INJ. 1μl
INSTRUMENT ID. SL6CC

0
0

RUN # 105 JAN/12/87 15:45:04
WORKFILE ID: C
WORKFILE NAME:

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.34	332516	D BV	0.135	9.563
1.72	54260	D VV	0.152	1.561
1.95	143457	D VB	0.115	4.126
2.47	268530	BV	0.120	7.723
2.93	89916	VB	0.155	2.560
6.26	2589201	SBB	0.319	74.467

TOTAL HEIGHT= 3477000
MUL FACTOR= 1.0000E+00

Page 1

Received: 12/12/86

SPECTRIX

REPORT

WORK ORDER # 86-12-001

RESULTS by Sample

SAMPLE ID CAV-DW09-001		SAMPLE # 04 FRACTIONS: A			
Date & Time Collected 12/12/86			Category _____		
AG	<10 ug/L	AS	13.0 ug/L	BE	5 ug/L
CD	<5 ug/L	CN SU	<100 ug/L	CR	31 ug/L
CU	21 ug/L	FE	30400 ug/L	HG	<0.2 ug/L
NI	<29 ug/L	PB	39.9 ug/L	SB	<44 ug/L
SE	<5 ug/L	TL	<10 ug/L	ZN	99 ug/L

CHECKLIST

SECTION II - Sample Data Packets

- Sample No: CAV-0W14-001
- Organic Analysis Data Sheets (OADS)
 - VOA
 - SV
 - Pest.
 - Tentatively Identified Compounds
 - VOA
 - SV
 - Raw Data
 - VOA chromatogram
 - VOA quantitation report (surrogates)
 - SV chromatogram
 - SV quantitation report (surrogates)
 - Pest. chromatogram and integrator print out
 - Inorganic Analysis Data Sheets (IADS)

004074

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-0W14-00 ,

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: B61205705
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *ymn*

CASE NO.: -----
QC REPORT NO.: -----
CONTRACT NO.: -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V05A
DILUTION FACTOR: .1.00

CAS #	COMPOUND	UG/L	
C010	CHLOROMETHANE	10	U
C015	BROMOMETHANE	10	U
C020	VINYL CHLORIDE	10	U
C025	CHLOROETHANE	10	U
C030	METHYLENE CHLORIDE	5	U
C035	ACETONE	12	
C040	CARBON DISULFIDE	5	U
C045	1, 1-DICHLOROETHENE	5	U
C050	1, 1-DICHLOROETHANE	5	U
C055	TRANS-1, 2-DICHLOROETHENE	5	U
C060	CHLOROFORM	5	U
C065	1, 2-DICHLOROETHANE	5	U
C110	2-BUTANONE	10	
C115	1, 1, 1-TRICHLOROETHANE	5	U
C120	CARBON TETRACHLORIDE	5	U
C125	VINYL ACETATE	10	
C130	BROMODICHLOROMETHANE	5	U
C140	1, 2-DICHLOROPROPANE	5	U
C145	TRANS-1, 3-DICHLOROPROPENE	5	U
C150	TRICHLOROETHENE	5	U
C155	DISROMOCHLOROMETHANE	5	U
C160	1, 1, 2-TRICHLOROETHANE	5	U
C165	BENZENE	5	U
C170	CIS-1, 3-DICHLOROPROPENE	5	U
C175	2-CHLOROETHYL VINYL ETHER	10	U
C180	BROMOFORM	5	U
C210	2-HEXANONE	10	
C205	4-METHYL-2-FENTANONE	10	U
C220	TETRACHLOROETHENE	5	U
C225	1, 1, 2-TETRACHLOROETHANE	5	U
C230	TOLUENE	9	U
C235	CHLOROBENZENE	5	U
C240	ETHYLBENZENE	10	
C245	STYRENE	5	U
C250	TOTAL XYLENES	37	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: CAV-DW14 - 001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205705
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *EPH*...

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86..

SEMICOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: ... 12/12/86.
 DATE ANALYZED: 01/12/87

DATAFILE: 4U12057C05

DILUTION FACTOR: 100.0...

CAS #	COMPOUND	UG/L
C315	PHENOL	1000 U
C325	BIS(2-CHLOROETHYL)ETHER	1000 U
C330	2-CHLOROPHENOL	1000 U
C335	1, 3-DICHLOROBENZENE	1000 U
C340	1, 4-DICHLOROBENZENE	1000 U
C345	BENZYL ALCOHOL	1000 U
C350	1, 2-DICHLOROBENZENE	1000 U
C355	2-METHYLPHENOL	1000 U
C360	BIS(2-CHLOROISOPROPYL)ETHER	1000 U
C365	4-METHYLPHENOL	1000 U
C370	N-NITROSODIPOXYLAMINE	1000 U
C375	HEXACHLOROETHANE	1000 U
C410	NITROBENZENE	1000 U
C415	ISOPHORONE	1000 U
C420	2-NITROPHENOL	1000 U
C425	2, 4-DIMETHYLPHENOL	1000 U
C430	BENZOIC ACID	5000 U
C435	BIS(2-CHLOROETHOXY)METHANE	1000 U
C440	2, 4-DICHLOROPHENOL	1000 U
C445	1, 2, 4-TRICHLOROBENZENE	1000 U
C450	NAPHTHALENE	8200
C455	4-CHLOROANILINE	1000 U
C460	HEXACHLOROBUTADIENE	1000 U
C465	P-CHLORO-M-CRESOL	1000 U
C470	2-METHYLNAPHTHALENE	350 J
C510	HEXACHLOROCYCLOPENTADIENE	1000 U
C515	2, 4, 6-TRICHLOROPHENOL	1000 U
C520	2, 4, 5-TRICHLOROPHENOL	5000 U
C525	2-CHLORONAPHTHALENE	1000 U
C530	2-NITROANILINE	5000 U
C535	DIMETHYL PHTHALATE	1000 U
C540	ACENAPHTHYLENE	1000 U
C545	3-NITROANILINE	5000 U
C550	ACENAPHTHENE	1000
C555	2, 4-DINITROPHENOL	5000 U
C560	4-NITROPHENOL	5000 U
C565	DIBENZOFURAN	600 J
C570	2, 4-DINITROTOLUENE	1000 U
C575	2, 6-DINITROTOLUENE	1000 U

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-DW14-00

EMI VOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C05

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	1000 U
C585	4-CHLOROPHENYL PHENYL ETHER	1000 U
C590	FLUORENE	610 J
C595	4-NITROANILINE	5000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	5000 U
C615	N-NITROSODIPHENYLAMINE	1000 U
C625	4-BROMOPHENYL PHENYL ETHER	1000 U
C630	HEXA CHLOROBENZENE	1000 U
C635	PENTACHLOROPHENOL	5000 U
C640	PHENANTHRENE	650 J
C645	ANTHRACENE	130 J
C650	DI-N-BUTYL PHTHALATE	1000 U
C655	FLUORANTHENE	360 J
C715	PYRENE	230 J
C720	BUTYL BENZYL PHTHALATE	1000 U
C725	3, 3'-DICHLOROBENZIDINE	2000 U
C730	BENZO(A)ANTHRACENE	1000 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	1000 U
C740	CHRYSENE	1000 U
C760	DI-N-OCTYL PHTHALATE	1000 U
C765	BENZO(B)FLUORANTHENE	1000 U
C770	BENZO(K)FLUORANTHENE	1000 U
C775	BENZO(A)PYRENE	1000 U
C780	INDENO(1, 2, 3-CD)PYRENE	1000 U
C785	DIBENZO(A, H)ANTHRACENE	1000 U
C790	BENZO(GHI)PERYLENE	1000 U

004077

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-05A

DC No: KEYSTONE
SMO No: CAV-OW14-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc:	Low	GPC Cleanup	Yes	No
Date Extracted:	12/18/86	Sep. Funnel Ext.	Yes	
Date Analyzed:	1/6/87	Cont. Liquid Ext.	Yes	
Dilution Factor:	1			
Percent Moisture(decanted):				

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOL-1016	0.50 U
11104-28-2	AROCLOL-1221	0.50 U
11141-16-5	AROCLOL-1232	0.50 U
53469-21-9	AROCLOL-1242	0.50 U
12672-29-6	AROCLOL-1248	0.50 U
11097-69-1	AROCLOL-1254	1.0 U
11096-82-5	AROCLOL-1260	1.0 U

004078

V_s = 500 = Volume of water extracted (ml)
V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000 /A.

2nd Column Inst. : 6000

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: CAV-0W14-00

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO. : -----

QC REPORT NO. : -----

ANALYST: CKT

DATAFILE: 4U12057V05A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.	00407	UG/	0

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: CAV-DW14

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: QU

DATAFILE: 4U12057C05

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMICOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
95-15-8	BENZO[B]THIOPHENE	746	904	0.450 J
				0
				0

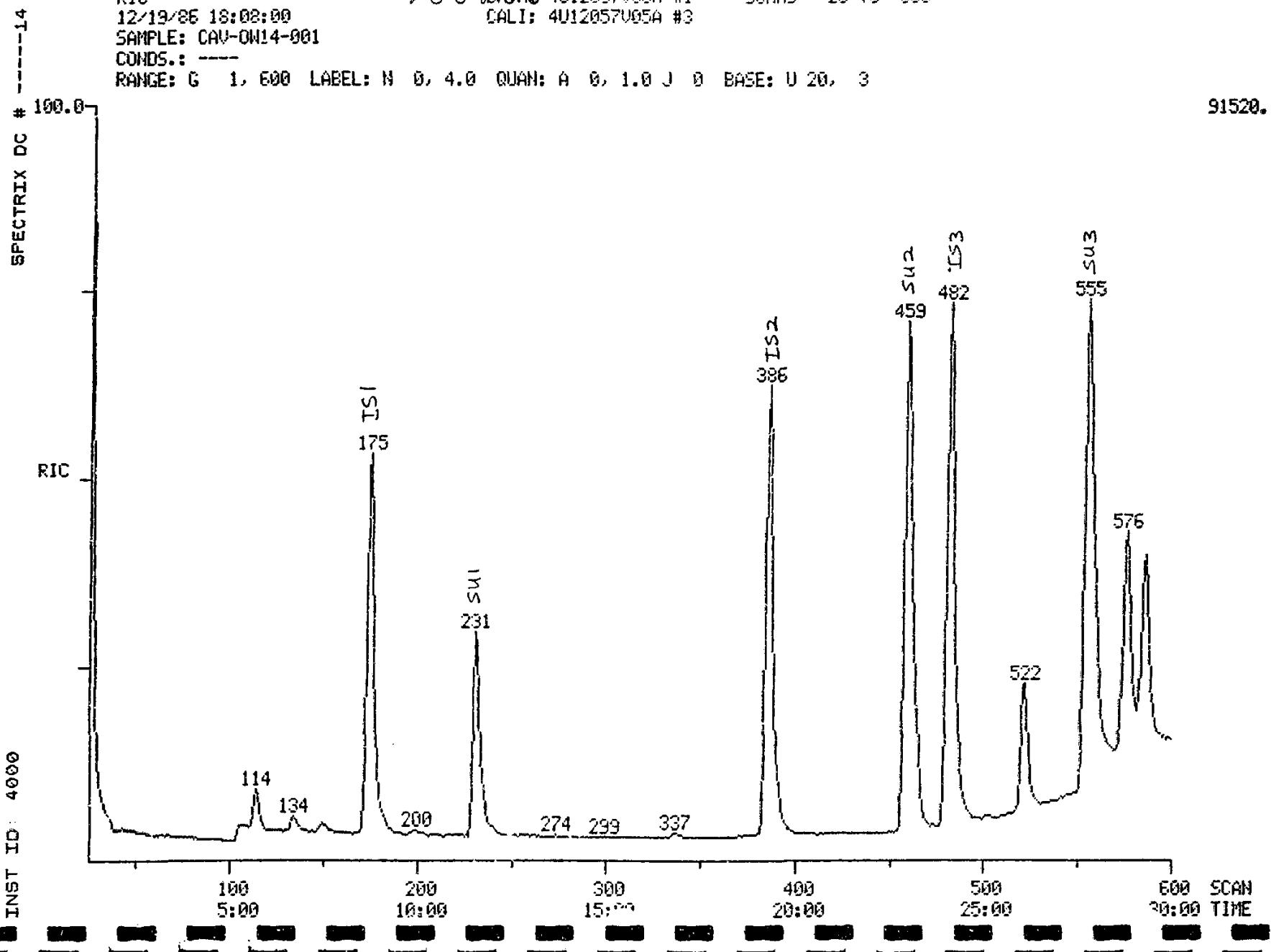
J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

RIC
12/19/86 18:02:00
SAMPLE: CAV-ON14-901
COND.: ----

L 8 0 TO 040 4U12057U05A #1 SCANS 25 TO 600
CALI: 4U12057U05A #3

RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

91520.



SPECTRIX DC # -----14

VOA QUANTITATION REPORT

AMPLE ID.	CAV-OW14-001		
FILENAME	4U12057V05A	INST	4000
CLIENT	KEYSTONE	ANALYST	CKT
DATE INJECTED	12/19/86 18:08:00	VERIFIED BY	<i>ypm</i>
CALIB STD DATE	12-14-86	CORR. FACTOR	1.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
174 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000 22317
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000 97764
482 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000 88271
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	84 %	1.328 27673
459 CS05	D8-TOLUENE	**SU2**	98	98 %	0.952 82657
555 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	91 %	1.151 54897
114 C030	METHYLENE CHLORIDE *		84	3 UG/L	0.655 2809
134 C035	ACETONE		43	12 UG/L	0.770 @009
463 C230	TOLUENE		92	9 UG/L	0.961 18210
521 C240	ETHYLBENZENE		106	10 UG/L	1.081 7931
576 C250	TOTAL XYLEMES		106	37 UG/L	1.195 34657

* BLANK CORRECTED ON DATA SHEET

INST ID: 4530

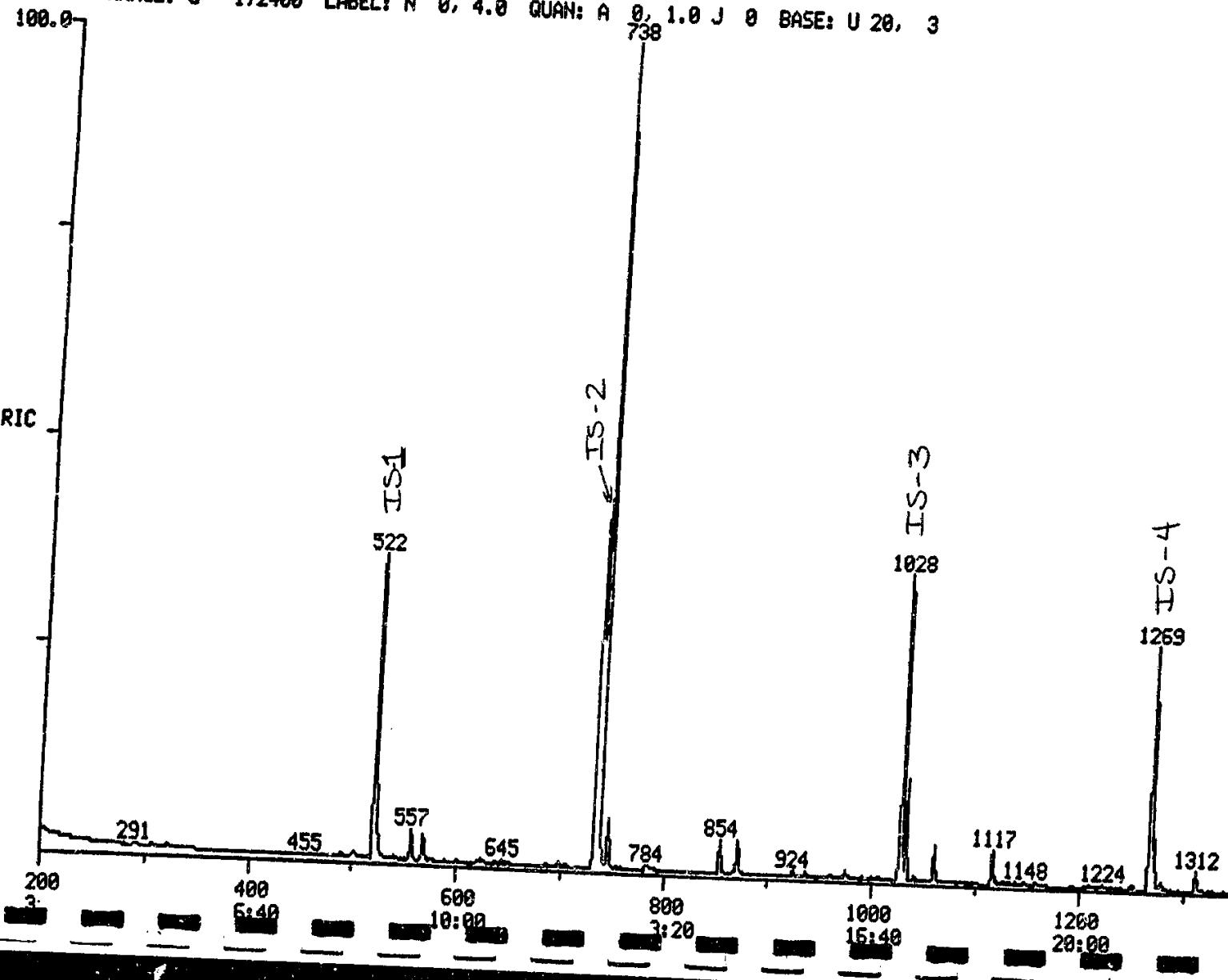
SPECTRUM DC # ---14

RIC
01/12/87 19:10:00
SAMPLE: CAV-OW14-001
COND.: ---
RANGE: G 1.2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

3 8 0 7 0 0

DATA: 4U12057C05 #1
CALI: 4U12057C05 #3

SCANS 260 TO 1350



004083

151552.

RIC
01/12/87 19:10:00
SAMPLE: CAV-0W14-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12057C05 #1
CALI: 4U12057C05 #3
SCANS 1350 TO 2400

21024.

SPECTRUM DC # -----14

100.0

RIC

INST ID: 4530

1453

1504

1408
23:20

1600
26:40

1800
36:00

2000
33:20

2200
36:40

2400
37:00

TS-5

TS-6

1928

2161 2242 2316

SCAN
TIME

004084

SPECTRIX DC # ----14

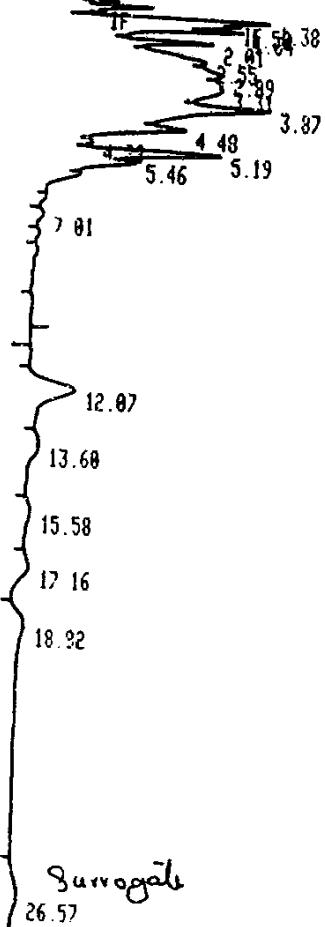
SEMOVOLATILE WATER REPORT

SAMPLE ID.	CAV-DW14-001		
FILENAME	4U12057C05	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/12/87 19:10:00	VERIFIED BY	<i>YPM</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	100.00

SCAN#	SEMOVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
521	CI30	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	4000 UG/L	1.000	2223%
735	CI40	NAPHTHALENE-D8 ** IS2 **	136	4000 UG/L	1.000	86621
1028	CI50	ACENAPHTHENE-D10 ** IS3 **	164	4000 UG/L	1.000	40326
1269	CI60	PHENANTHRENE-D10 ** IS4 **	188	4000 UG/L	1.000	5150%
1709	CI70	CHRYSENE-D12 ** IS5 **	240	4000 UG/L	1.000	002946%
1928	CI75	PERYLENE-D12 ** IS6 **	264	4000 UG/L	1.000	01411%
738	C450	NAPHTHALENE	128	8197 UG/L	1.004	17387%
854	C470	2-METHYLNAPHTHALENE	142	354 UG/L	1.162	5086
1033	C550	ACENAPHTHENE	153	1046 UG/L	1.005	9940
1060	C565	DIBENZOFURAN	168	596 UG/L	1.031	7510
1117	C590	FLUORENE	166	611 UG/L	1.087	6512
1272	C640	PHENANTHRENE	178	652 UG/L	1.002	815%
1279	C645	ANTHRACENE	178	129 UG/L	1.008	163%
1469	C655	FLUORANTHENE	202	360 UG/L	1.158	393%
504	C715	PYRENE	202	226 UG/L	0.880	260%

* SURROGATES WERE DILUTED OUT

STK-L-86-12-057-05A 1:2 NA.



DC#
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAY-QW14-001
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOSORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1.44
INSTRUMENT ID. 6000

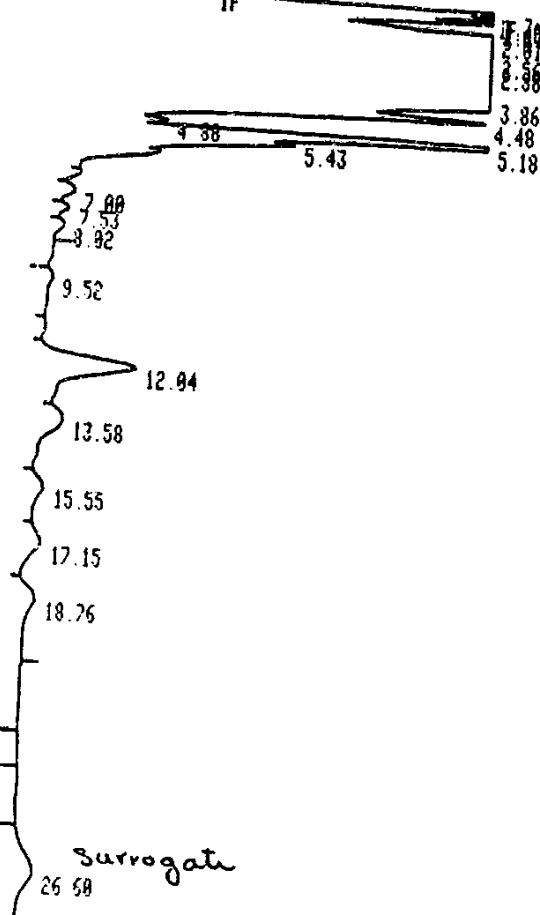
0
0

RUN # 22 JAN/06/87 23:52:10
WORKFILE ID: C
WORKFILE NAME:

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.38	47725	BY	0.119	10.149
1.50	40237	D VV	0.111	8.556
1.64	41289	D VV	0.163	8.780
2.01	34568	VV	0.120	7.351
2.55	33700	VV	0.372	7.166
2.89	38928	VV	0.355	6.278
3.31	39809	VV	0.646	8.465
3.87	53562	VV	0.435	11.390
4.48	33009	VV	0.284	7.019
4.88	9754	VV	0.193	2.074
5.19	43523	VV	0.227	9.255
5.46	22121	VV	0.264	4.704
7.01	1424	PP	0.291	0.303
12.07	12724	VV	0.735	2.706
13.60	4317	VV	1.223	0.918
15.58	3337	VV	1.132	0.710
17.16	3955	VP	0.947	0.841
18.82	3429	PV	2.981	0.729
26.57	2845	VP	1.955	0.605

STK# 44 56-12-057-05A

N.A.



DC# _____
SPECTRIX HOUSTON
CASE NO. Keystone

SAMPLE NO. CAV-0W14-001
PESTICIDE FRACTION
2ND COLUMN CONFIRMA

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187°C
VOLUME INJ. 1.46
INSTRUMENT ID. 6900

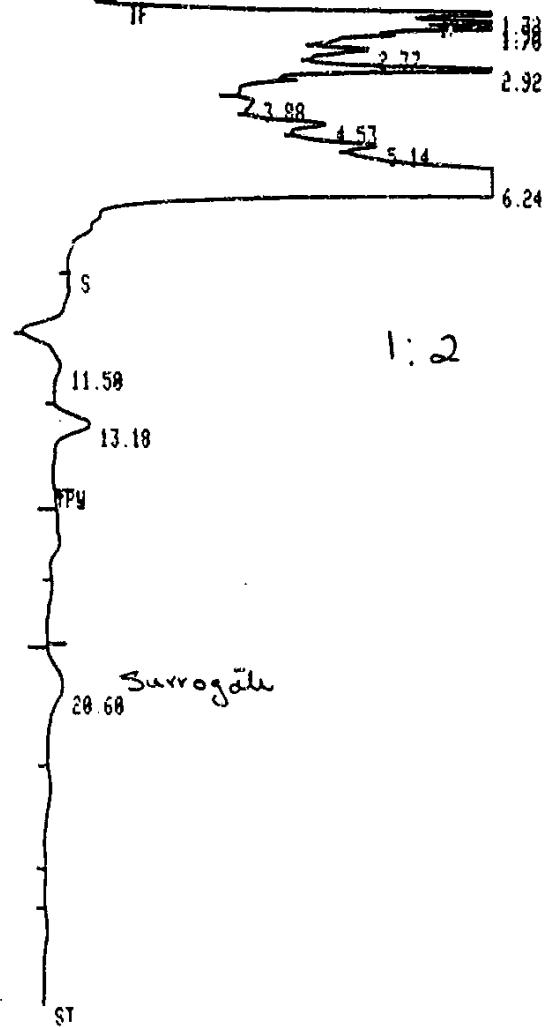
4
0
0

RUN # 24 JAN/07/87 01:04:42
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.36	119160	PV	0.112	8.766
1.49	96196	D VV	0.118	7.076
1.64	94320	D VV	0.131	6.938
2.01	132167	VV	0.230	9.722
2.56	171808	VV	0.390	12.638
2.83	207840	VV	0.395	15.289
3.86	189434	VV	0.395	13.935
4.43	94725	VV	0.239	6.968
4.88	14911	VV	0.181	1.097
5.18	125500	VV	0.216	9.232
5.43	59780	D VV	0.231	3.735
7.00	2946	PP	0.301	0.209
7.92	2411	PV	0.228	0.177
9.52	2060	VB	0.260	0.152
12.04	1542	BY	0.619	0.113
13.58	25218	VV	0.693	1.855
15.55	7495	VV	0.994	0.551
17.15	4033	VV	0.891	0.297
18.26	4618	VP	0.897	0.340
26.60	3934	PB	0.866	0.282
34.67	4913	BB	1.396	0.361
	3598	I BH	1.513	0.265

STANLIF 86-12-053-05A 1:2 NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-6W14.00
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm c
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 60°C
VOLUME INJ. 1uL
INSTRUMENT ID. 600

RUN # 197 JAN/12/87 16:47:37
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.32	104703	D BV	0.121	5.018
1.53	50736	D VV	0.147	2.431
1.76	39326	D VV	0.172	1.885
2.37	15754	VV	0.297	0.755
2.92	47836	VB	0.176	2.292
3.88	1974	PV	0.302	0.895
4.53	11162	VV	0.343	0.535
5.14	17829	VH	0.373	0.854
6.24	1785215	SHB	0.281	85.550
11.58	3959	PV	1.306	0.190
13.18	6142	VB	0.910	0.294
20.68	2107	BV	1.373	0.161

Page 1
Received: 12/12/86

SPECTRIX REPORT

Work Order # 86-12-057

68070 Results by Sample

SAMPLE ID <u>CAV-DW14-001</u>		SAMPLE # <u>05</u> FRACTIONS: A										
		Date & Time Collected <u>12/12/86</u>										
AG	<10	AS	83.0	BE	6	CD	<5	CN	SU	<100	CR	68
CU	14	FE	39500	HG	<0.2	NJ	<29	PB	28.6	SB	<44	
SE	<5	TL	<10	ZN	121							

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-MW04-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

004090

SPECTRIX DJ # ----- 8

INST ID 4000

SAMPLE NUMBER: SCK-MW04-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205706
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *SPH*

CASE NO.: -----
QC REPORT NO.: -----
CONTRACT NO.: -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V06
DILUTION FACTOR: 1:20

CAS #	COMPOUND	UG/L	1
C010	CHLOROMETHANE	200	U
C015	BROMOMETHANE	200	U
C020	VINYL CHLORIDE	200	U
C025	CHLOROETHANE	200	U
C030	METHYLENE CHLORIDE	180	
C035	ACETONE	200	U
C040	CARBON DISULFIDE	100	U
C045	1, 1-DICHLOROETHENE	100	U
C050	1, 1-DICHLOROETHANE	100	U
C055	TRANS-1, 2-DICHLOROETHENE	100	U
C060	CHLOROFORM	100	U
C065	1, 2-DICHLOROETHANE	100	U
C110	2-BUTANONE	200	U
C115	1, 1, 1-TRICHLOROETHANE	100	U
C120	CARBON TETRACHLORIDE	100	U
C125	VINYL ACETATE	200	U
C130	BROMODICHLOROMETHANE	100	U
C140	1, 2-DICHLOROPROPANE	100	U
C145	TRANS-1, 3-DICHLOROPROPENE	100	U
C150	TRICHLOROETHENE	100	U
C155	DIBROMOCHLOROMETHANE	100	U
C160	1, 1, 2-TRICHLOROETHANE	100	U
C165	BENZENE	50	J
C170	CIS-1, 3-DICHLOROPROPENE	100	U
C175	2-CHLOROETHYL VINYL ETHER	200	U
C180	BROMOFORM	100	U
C210	2-HEXANONE	200	U
C205	4-METHYL-2-PENTANONE	200	U
C220	TETRACHLOROETHENE	100	U
C225	1, 1, 2, 2-TETRACHLOROETHANE	100	U
C230	TOLUENE	100	U
C235	CHLOROBENZENE	100	U
C240	ETHYLBENZENE	100	U
C245	STYRENE	100	U
C250	TOTAL XYLEMES	100	U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # --- 8
SAMPLE NUMBER: SCK-MW04 - 001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205706
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *YJZ*

CASE NO. : ---
 QC REPORT NO. : ---
 CONTRACT NO. : ---
 DATE SAMPLE RECEIVED: 12/12/86.

SEMOVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/18/86.
 DATE ANALYZED: 01/13/87

DATAFILE: 4U12057C06A

DILUTION FACTOR: ...5.0...

CAS #	COMPOUND	UG/L	
C315	PHENOL	50 U	
C325	BIS(2-CHLOROETHYL)ETHER	50 U	
C330	2-CHLOROPHENOL	50 U	
C339	1,3-DICHLOROBENZENE	50 U	
C340	1,4-DICHLOROBENZENE	50 U	
C345	BENZYL ALCOHOL	50 U	
C350	1,2-DICHLOROBENZENE	50 U	
C355	2-METHYLPHENOL	50 U	
C360	BIS(2-CHLOROISOPROPYL)ETHER	50 U	
C365	4-METHYLPHENOL	50 U	
C370	N-NITROSODIISOPROPYLAMINE	50 U	
C375	HEXACHLOROETHANE	50 U	
C410	NITROBENZENE	50 U	
C415	ISOPHORONE	50 U	
C420	2-NITROPHENOL	50 U	
C425	2,4-DIMETHYLPHENOL	50 U	
C430	BENZOIC ACID	250 U	
C435	BIS(2-CHLOROETHOXY)METHANE	50 U	
C440	2,4-DICHLOROPHENOL	50 U	
C445	1,2,4-TRICHLOROBENZENE	50 U	
C450	NAPHTHALENE	540	
C455	4-CHLOROANILINE	50 U	
C460	HEXACHLOROBUTADIENE	50 U	
C465	P-CHLORO-M-CRESO	50 U	
C470	2-METHYLNAPHTHALENE	7 U	
C510	HEXACHLOROCYCLOPENTADIENE	50 U	
C515	2,4,6-TRICHLOROPHENOL	50 U	
C520	2,4,5-TRICHLOROPHENOL	250 U	
C525	2-CHLORONAPHTHALENE	50 U	
C530	2-NITROANILINE	250 U	
C535	DIMETHYL PHTHALATE	50 U	
C540	ACENAPHTHYLENE	50 U	
C545	3-NITROANILINE	250 U	
C550	ACENAPHTHENE	150	
C555	2,4-DINITROPHENOL	250 U	
C560	4-NITROPHENOL	250 U	
C565	DIBENZOFURAN	77	
C570	2,4-DINITROTOLUENE	50 U	
C575	2,6-DINITROTOLUENE	50 U	

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-MW04 -00

EMILOVATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C06A

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	50 U
C585	4-CHLOROPHENYL PHENYL ETHER	50 U
C590	FLUORENE	84
C595	4-NITROANILINE	250 U
C610	4, 6-DINITRO-2-METHYLPHENOL	250 U
C615	N-NITROSODIPHENYLAMINE	50 U
C625	4-BROMOPHENYL PHENYL ETHER	50 U
C630	HEXACHLOROBENZENE	50 U
C635	PENTACHLOROPHENOL	250 U
C640	PHENANTHRENE	21 J
C645	ANTHRACENE	8 J
C650	DI-N-BUTYL PHTHALATE	50 U
C655	FLUORANTHENE	50 U
C715	PYRENE	50 U
C720	BUTYL BENZYL PHTHALATE	50 U
C725	3, 3'-DICHLOROBENZIDINE	100 U
C730	BENZO(A)ANTHRACENE	50 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	50 U
C740	CHRYSENE	50 U
C760	DI-N-OCTYL PHTHALATE	50 U
C765	BENZO(B)FLUORANTHENE	50 U
C770	BENZO(K)FLUORANTHENE	50 U
C775	BENZO(A)PYRENE	50 U
C780	INDENO(1, 2, 3-CD)PYRENE	50 U
C785	DIBENZO(A, H)ANTHRACENE	50 U
C790	BENZO(ghi)PERYLENE	50 U

004003

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-06A

DC No: KEYSTONE
SMO No: SCK-MW04-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low

GPC Cleanup Yes No

Date Extracted: 12/18/86

Sep. Funnel Ext. Yes

Date Analyzed: 1/7/87

Cont. Liquid Ext. Yes

Dilution Factor: 1

Percent Moisture(decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-33-4	ENDRIN ALDEHYDE	0.10 U
1031-07-6	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOR-1016	0.50 U
11104-28-2	AROCLOR-1221	0.50 U
11141-16-5	AROCLOR-1232	0.50 U
53469-21-9	AROCLOR-1242	0.50 U
12672-29-6	AROCLOR-1248	0.50 U
11097-69-1	AROCLOR-1254	1.0 U
11096-82-5	AROCLOR-1260	1.0 U

004094

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 N/A.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-MW04-00

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO. : -----

QC REPORT NO. : -----

ANALYST: NOT

DATAFILE: 4U12057V06

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
		0	UG/	0
		0	0	0

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: SCK-MW04 - 001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12057C06A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	305		0UG/L
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	563	957	038 J
95-13-6	1H-INDENE	574	933	4240 J
95-15-8	BENZO[B]THIOPHENE	750	917	068 J
2471-83-2	1H-INDENE, 1-ETHYLIDENE-	874	900	27 J
86-74-8	9H-CARBAZOLE	1315	885	36 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

RIC
12/19/86 14:22:00
SAMPLE: SCK-MN04-001
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

L 6 0 DATA: 04U12057U06 #1
CALI: 4U12057U06 #3
SCANS 25 TO 600

56192.

SPECTRIX DC # -----14

RIC

INST ID: 4000

100.0

50.0

10.0

1.0

0.1

0.0

TS1

122

155

100

5:00

200

10:00

300

15:00

400

20:00

500

25:00

600
29:00
SCAN
TIME

TS1

233

201

285

311

337

366

TS2

386

TS3

459

482

TS3

556

522

586

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

AMPLE ID.	SCK-MW04-001		
FILENAME	4U12057V06	INST	4000
CLIENT	KEYSTONE	ANALYST	NOT
DATE INJECTED	12/19/86 14:22:00	VERIFIED BY	NOT
CALIB STD DATE	12-14-86	CORR. FACTOR	20.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
178 CI01	BROMOCHLOROMETHANE	**IS1**	128	1000 UG/L	1.000
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	1000 UG/L	1.000
482 CI20	D5-CHLOROBENZENE	**IS3**	117	1000 UG/L	1.000
233 CS15	D4-1, 2-CHLOROETHANE	**SU1**	65	86 %	1.309
459 CS05	D8-TOLLENE	**SU2**	98	100 %	0.952
555 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	92 %	1.151
122 C030	METHYLENE CHLORIDE*		84	279 UG/L	0.685
337 C165	BENZENE		78	50 UG/L	0.873

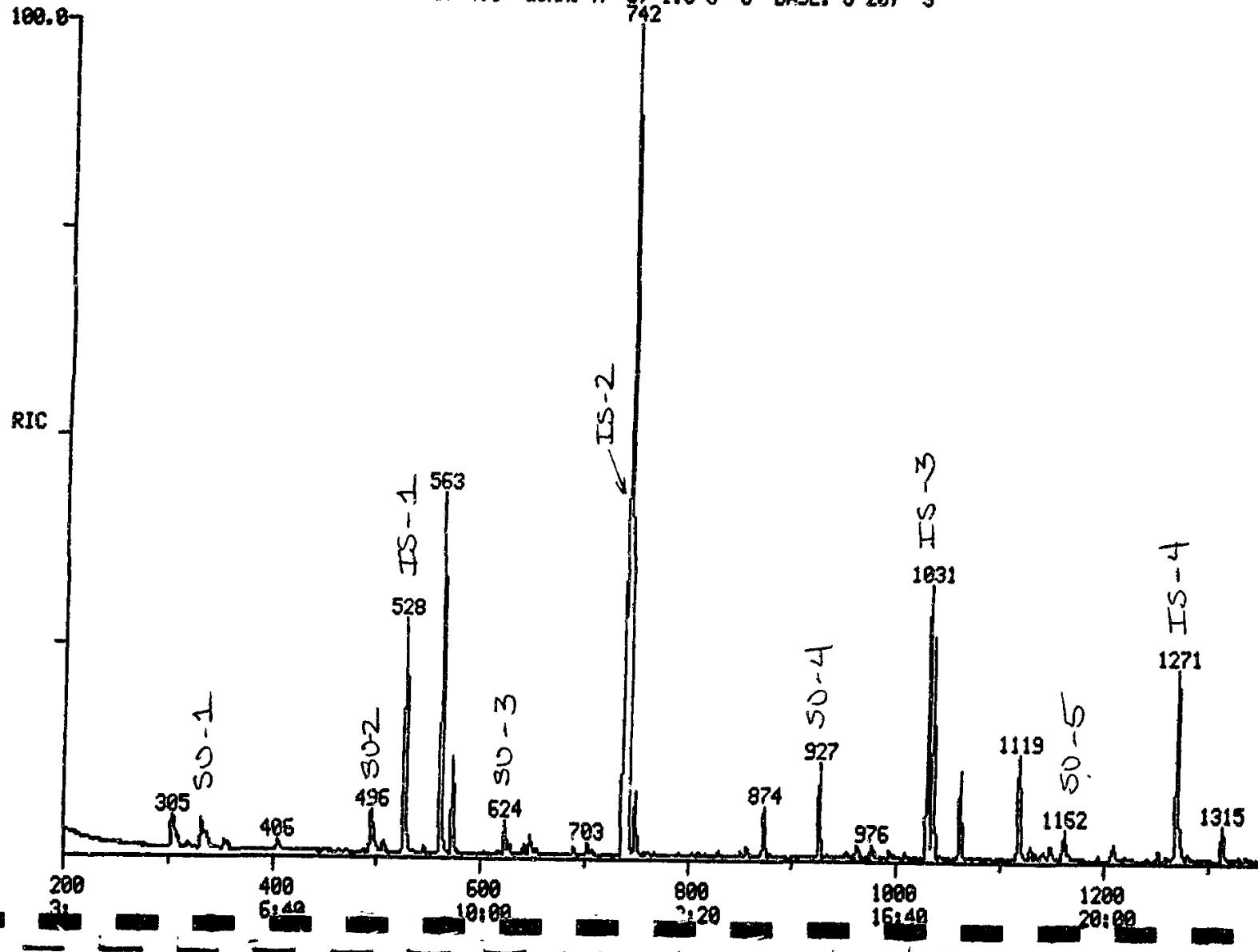
* BLANK CORRECTED ON DATA SHEET

INST-ID: 4530

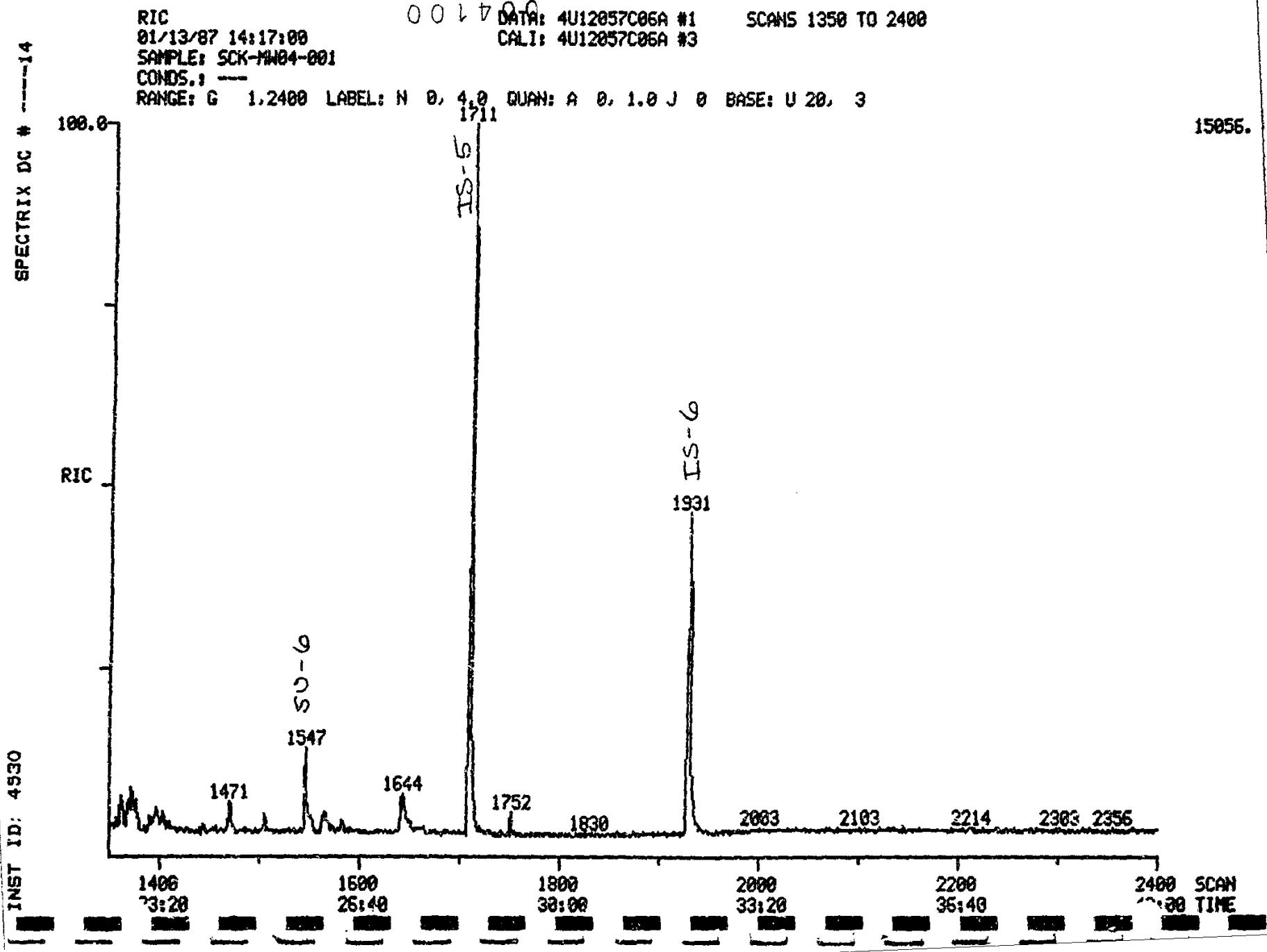
SPECTRIX DC # ---14

RIC
01/13/87 14:17:00
SAMPLE: SCK-MM04-001
COND.: ---
RANGE: G 1.2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

660 700
DATA: 4U12057C06A #1
CALI: 4U12057C06A #3
SCANS 200 TO 1350



151808.



SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	SCK-MW04-001		
FILENAME	4U12057C06A	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/13/87 14:17:00	VERIFIED BY	<i>MTH</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	5.00

SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
528	CI30	1,4-DICHLOROBENZEND-D4 ** IS1 **	132	200 UG/L	1.000	1877
738	CI40	NAPHTHALENE-D8 ** IS2 **	136	200 UG/L	1.000	7005
1031	CI50	ACENAPHTHENE-D10 ** IS3 **	164	200 UG/L	1.000	3564
1271	CI60	PHENANTHRENE-D10 ** IS4 **	188	200 UG/L	1.000	4325
1711	CI70	CHRYSENE-D12 ** IS5 **	240	200 UG/L	1.000	2052
1931	CI75	PERYLENE-D12 ** IS6 **	264	200 UG/L	1.000	1316
333	CS50	2-FLUOROPHENOL ** SU1 **	112	60 %	0.631	750
495	CS45	PHENOL-D5 ** SU2 **	99	63 %	0.937	1167
1162	CS55	2,4,6-TRIBROMOPHENOL ** SU5 **	330	64 %	1.127	131
624	CS20	NITROBENZENE-D5 ** SU3 **	82	79 %	0.846	642
927	CS25	2-FLUOROBIPHENYL ** SU4 **	172	75 %	0.899	426
1547	CS30	TERPHENYL-D14 ** SU6 **	244	40 %	0.904	218
742	C450	NAPHTHALENE	128	543 UG/L	1.005	18919
857	C470	2-METHYLNAPHTHALENE	142	7 UG/L	1.161	155
036	C550	ACENAPHTHENE	153	154 UG/L	1.005	2437
1062	C565	DIBENZOFURAN	168	77 UG/L	1.030	1703
1119	C590	FLUORENE	166	84 UG/L	1.085	1574
1274	C640	PHENANTHRENE	178	21 UG/L	1.002	436
1282	C645	ANTHRACENE	178	8 UG/L	1.009	177

INST ID: 4530

SPECTRUM DC # ---19

1004
SAMPLE

LIBRARY SEARCH
01/13/87 14:17:00 + 5:05
SAMPLE: SCK-MN04-001
COND.: --
ENHANCED (5 15B 2N 0T)

004102
DATA: 4U12057C06A # 305
CALI: 4U12057C06A # 3
BASE M/Z: 91
RIC: 4751.

C8.H18
M WT 1004
B PK 91
RANK 1
1725
PUR 694

BENZENE, ETHYL-

C8.H10
M WT 1004
B PK 91
RANK 2
1739
PUR 682

1,3-CYCLOPENTADIENE, 5-(1-METHYLETHYLIDENE)-

C8.H10
M WT 1004
B PK 91
RANK 3
1724
PUR 664

BENZENE, 1,2-DIMETHYL-

M/Z

30 40 50 60 70 80 90 100

START # 86-12-057-06A

NA.

2.48
2.52
3.86
4.89

2.04

DC#
SPECTRIX HOUSTON
CASE NO. Keystone

SAMPLE NO. SCK-MW04-O
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 μL
INSTRUMENT ID. 6500

4
0
0

Surrogate
26.58

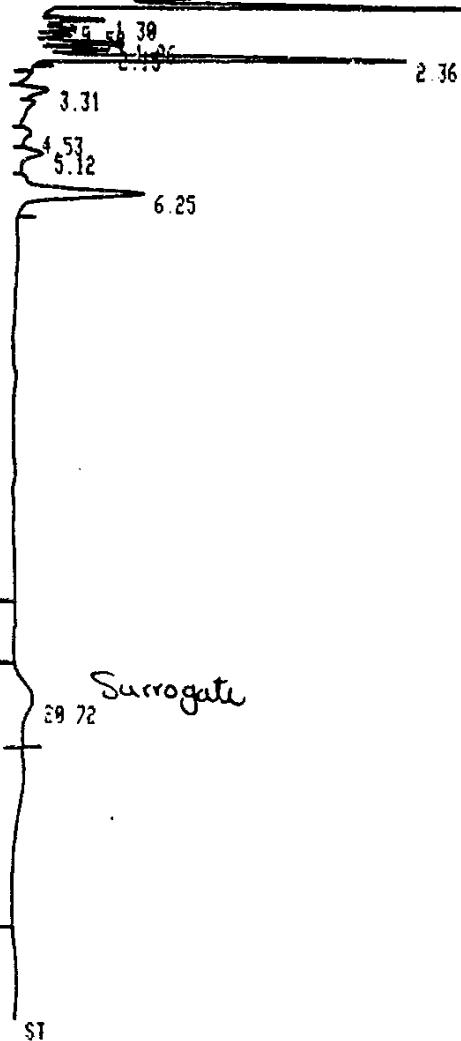
ST

RUN # 26
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%
RT HEIGHT TYPE AR/HT HEIGHT%
1.36 6677 BV 0.086 3.202
1.49 11376 D VV 0.115 5.455
1.64 46023 D VV 0.104 22.069
2.00 109196 D VB 0.105 52.362
2.48 9974 BP 0.167 4.783
2.86 3584 PV 0.151 1.719
3.12 2988 D VB 0.141 1.433
3.86 11264 BV 0.210 5.401
4.89 3471 VV 0.227 1.664
26.58 3988 PP 1.450 1.912

WFL: 86-12-057-06A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. SCK-MW04-C
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x 6mm C
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 603°C
VOLUME INJ. Full
INSTRUMENT ID. 0600

O

RUN # 108 JAN/12/87 17:18:53
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.38	6608	D BY	0.086	6.243
1.58	3220	D VV	0.090	3.042
1.76	3294	D VV	0.088	3.112
1.96	10316	D VV	0.100	9.746
2.15	8625	D VV	0.107	8.148
2.36	47364	V8	0.118	44.745
3.31	3623	PV	0.223	3.423
4.53	1665	VV	0.334	1.573
5.12	3102	VV	0.286	2.931
6.25	16050	VB	0.311	15.162
29.72	1987	PR	0.968	1.877

Page 1
Received: 12/12/86

SPECTRIX REPORT
S O L A R Results by Sample

Work Order # 86-12-05/

SAMPLE ID <u>SCK-MW04-001</u>		SAMPLE # 06 FRACTIONS: A									
		Date & Time Collected <u>12/12/86</u>								Category	
AG	<10	AS	25.3	BE	4	CD	<5	CN	SU	<100	CR
	ug/L		ug/L		ug/L		ug/L		ug/L		ug/L
CU	<10	FE	7510	HG	<0.2	NI	<29	PB	36.4	SB	<44
	ug/L		ug/L		ug/L		ug/L		ug/L		ug/L
SE	<5	TL	<10	ZN	61						
	ug/L		ug/L		ug/L						

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-MW10-001

Organic Analysis Data Sheets (OADS)

- VOA
- SV
- Pest.

Tentatively Identified Compounds

- VOA
- SV

Raw Data

- VOA chromatogram
- VOA quantitation report (surrogates)
- SV chromatogram
- SV quantitation report (surrogates)
- Pest. chromatogram and Integrator print out

Inorganic Analysis Data Sheets (IADS)

004106

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: SCK-MW10-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205707
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *John*

CASE NO.: -----
QC REPORT NO.: -----
CONTRACT NO.: -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V07

DILUTION FACTOR: 1:5

CAS #	COMPOUND	UG/L	004107
C010	CHLOROMETHANE	50 U	
C015	BROMOMETHANE	50 U	
C020	VINYL CHLORIDE	50 U	
C025	CHLOROETHANE	50 U	
C030	METHYLENE CHLORIDE	25 U	
C035	ACETONE	50 U	
C040	CARBON DISULFIDE	25 U	
C045	1, 1-DICHLOROETHENE	25 U	
C050	1, 1-DICHLOROETHANE	25 U	
C055	TRANS-1, 2-DICHLOROETHENE	25 U	
C060	CHLOROFORM	25 U	
C065	1, 2-DICHLOROETHANE	25 U	
C110	2-BUTANONE	50 U	
C115	1, 1, 1-TRICHLOROETHANE	25 U	
C120	CARBON TETRACHLORIDE	25 U	
C125	VINYL ACETATE	50 U	
C130	BROMODICHLOROMETHANE	25 U	
C140	1, 2-DICHLOROPROPANE	25 U	
C145	TRANS-1, 3-DICHLOROPROPENE	25 U	
C150	TRICHLOROETHENE	25 U	
C155	DIBROMOCHLOROMETHANE	25 U	
C160	1, 1, 2-TRICHLOROETHANE	25 U	
C165	BENZENE	8 U	
C170	CIS-1, 3-DICHLOROPROPENE	25 U	
C175	2-CHLOROETHYL VINYL ETHER	50 U	
C180	BROMOFORM	25 U	
C210	2-HEXANONE	50 U	
C205	4-METHYL-2-PENTANONE	50 U	
C220	TETRACHLOROETHENE	25 U	
C225	1, 1, 2, 2-TETRACHLOROETHANE	25 U	
C230	TOLUENE	25 U	
C235	CHLOROBENZENE	25 U	
C240	ETHYLBENZENE	25 U	
C245	STYRENE	25 U	
C250	TOTAL XYLEMES	11 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # --- 8
SAMPLE NUMBER: SCK-MW10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205707
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *YPM*

CASE NO.: ---
QC REPORT NO.: ---
CONTRACT NO.: ---
DATE SAMPLE RECEIVED: 12/12/86..

SEMOVOLATILES

CONCENTRATION: LOW
DATE EXTRACTED: 12/18/86.
DATE ANALYZED: 01/13/87

DATAFILE: 4U12057C07A
DILUTION FACTOR: 20.0....

CAS #	COMPOUND	UG/L	
C315	PHENOL	200 U	
C325	BIS(2-CHLOROETHYL)ETHER	200 U	00
C330	2-CHLOROPHENOL	200 U	0
C335	1, 3-DICHLOROBENZENE	200 U	1
C340	1, 4-DICHLOROBENZENE	200 U	0
C345	BENZYL ALCOHOL	200 U	4
C350	1, 2-DICHLOROBENZENE	200 U	0
C355	2-METHYLPHENOL	200 U	0
C360	BIS(2-CHLOROISOPROPYL)ETHER	200 U	
C365	4-METHYLPHENOL	200 U	
C370	N-NITROSODIPROPYLAMINE	200 U	
C375	HEXAChLOROETHANE	200 U	
C410	NITROBENZENE	200 U	
C415	ISOPHORONE	200 U	
C420	2-NITROPHENOL	200 U	
C425	2, 4-DIMETHYLPHENOL	200 U	
C430	BENZOIC ACID	1000 U	
C435	BIS(2-CHLOROETHOXY)METHANE	200 U	
C440	2, 4-DICHLOROPHENOL	200 U	
C445	1, 2, 4-TRICHLOROBENZENE	200 U	
C450	NAPHTHALENE	3400	
C455	4-CHLOROANILINE	200 U	
C460	HEXAChLOROBUTADIENE	200 U	
C465	P-CHLORO-M-CRESOL	200 U	
C470	2-METHYLNAPHTHALENE	160 U	
C510	HEXAChLOROCYCLOPENTADIENE	200 U	
C515	2, 4, 6-TRICHLOROPHENOL	200 U	
C520	2, 4, 5-TRICHLOROPHENOL	1000 U	
C525	2-CHLORONAPHTHALENE	200 U	
C530	2-NITROANILINE	1000 U	
C535	DIMETHYL PHTHALATE	200 U	
C540	ACENAPHTHYLENE	200 U	
C545	3-NITROANILINE	1000 U	
C550	ACENAPHTHENE	680	
C555	2, 4-DINITROPHENOL	1000 U	
C560	4-NITROPHENOL	1000 U	
C565	DIBENZOFURAN	340	
C570	2, 4-DINITROTOLUENE	200 U	
C575	2, 6-DINITROTOLUENE	200 U	

SPECTRIX DC # --- 8

SAMPLE NUMBER: SCK-MW10-001

EMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C07A

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	200 U
C585	4-CHLOROPHENYL PHENYL ETHER	200 U
C590	FLUORENE	210
C595	4-NITROANILINE	1000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	1000 U
C615	N-NITROSODIPHENYLAMINE	200 U
C625	4-BROMOPHENYL PHENYL ETHER	200 U
C630	HEXACHLOROBENZENE	200 U
C635	PENTACHLOROPHENOL	1000 U
C640	PHENANTHRENE	410
C645	ANTHRACENE	50 J
C650	DI-N-BUTYL PHTHALATE	200 U
C655	FLUORANTHENE	56 J
C715	PYRENE	41 J
C720	BUTYL BENZYL PHTHALATE	200 U
C725	3, 3'-DICHLOROBENZIDINE	400 U
C730	BENZO(A)ANTHRACENE	200 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	200 U
C740	CHRYSENE	200 U
C760	DI-N-OCTYL PHTHALATE	200 U
C765	BENZO(B)FLUORANTHENE	200 U
C770	BENZO(K)FLUORANTHENE	200 U
C775	BENZO(A)PYRENE	200 U
C780	INDENO(1, 2, 3-CD)PYRENE	200 U
C785	DIBENZO(A, H)ANTHRACENE	200 U
C790	BENZO(GHI)PERYLENE	200 U

004109

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-07A

DC No: KEYSTONE
SMO No: SCK-MW10-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc:	Low	GPC Cleanup	Yes	No
Date Extracted:	12/18/86	Sep. Funnel Ext.	Yes	/
Date Analyzed:	1/7/87	Cont. Liquid Ext.	Yes	/
Dilution Factor:	1			

Percent Moisture(decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCOLOR-1016	0.50 U
11104-28-2	AROCOLOR-1221	0.50 U
11141-16-5	AROCOLOR-1232	0.50 U
53469-21-9	AROCOLOR-1242	0.50 U
12672-29-6	AROCOLOR-1248	0.50 U
11097-69-1	AROCOLOR-1254	1.0 U
11096-82-5	AROCOLOR-1260	1.0 U

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-MW10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: NOT

DATAFILE: 4U12057V07

6. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
-------	-------------------------	-------	--------	--------

0041
UG/

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: SCK-MW10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12057C07A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	567	956	310 J
95-15-8	BENZO[B]THIOPHENE	752	914	250 J
2471-83-2	1H-INDENE, 1-ETHYLIDENE-	875	903	280 J
92-52-4	1,1'-BIPHENYL	940	929	220 J
81-40-8	NAPHTHALENE, 2,3-DIMETHYL-	977	934	120 J
86-74-8	9H-CARBAZOLE	1316	909	82 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

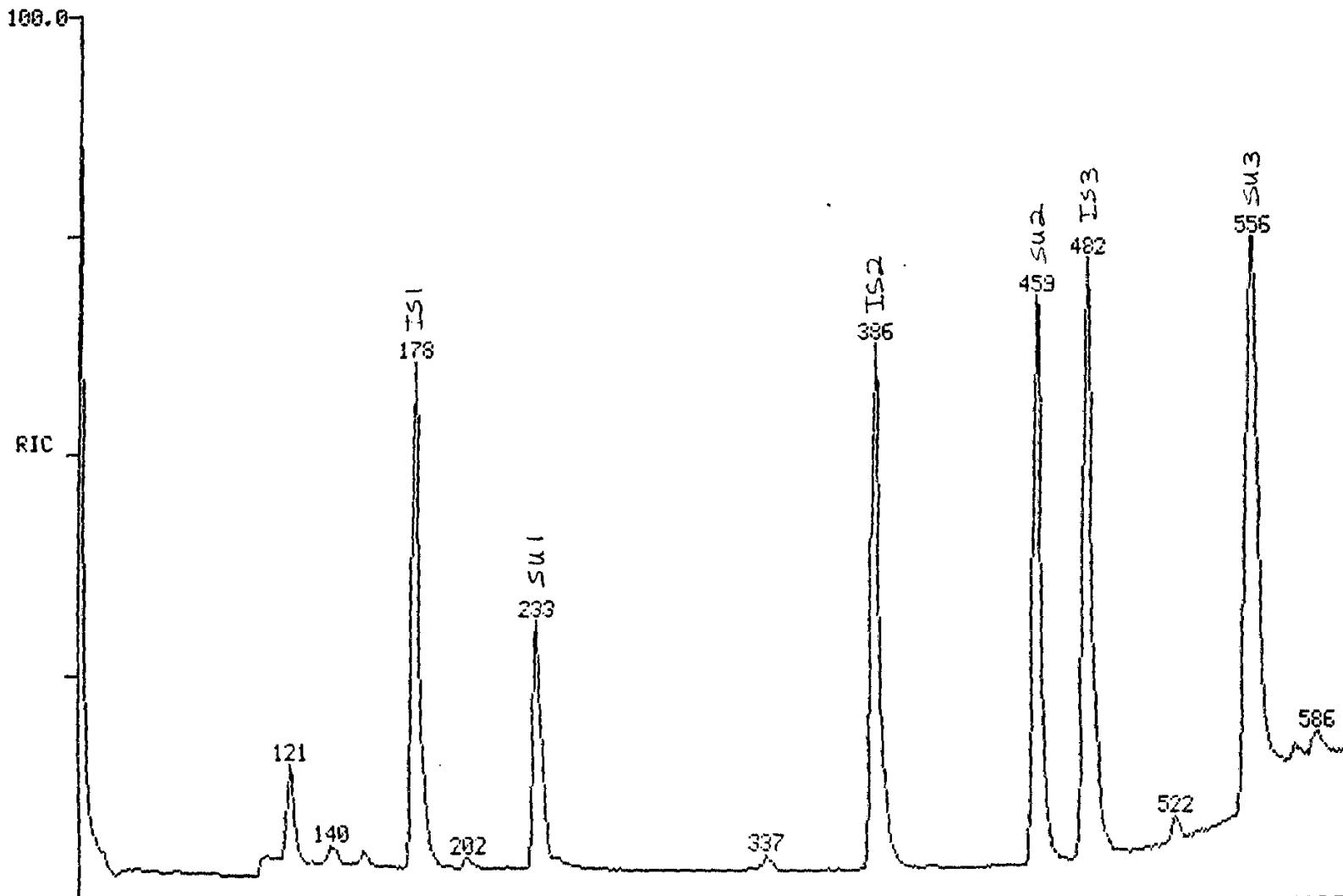
RIC
12/19/86 15:18:00
SAMPLE: SCX-MW10-001
CONDNS.: ----

DATA: 4U12057V07 #1
CALI: 4U12057V07 #3
SCANS 25 TO 600
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

84224.

SPECTRIX DC # -----14

INST ID: 4000



600 SCAN
30:00 TIME

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

AMPLE ID.	SCK-MW10-001		
FILENAME	4U12057V07	INST	4000
CLIENT	KEYSTONE	ANALYST	NOT
DATE INJECTED	12/19/86 15:18:00	VERIFIED BY	<i>Not</i>
CALIB STD DATE	12-14-86	CORR. FACTOR	5.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
178 CI01	BROMOCHLOROMETHANE	**IS1**	128	250 UG/L	1.000
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	250 UG/L	1.000
482 CI20	D5-CHLOROBENZENE	**IS3**	117	250 UG/L	1.000
233 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	84 %	1.309
459 CS05	D8-TOLUENE	**SU2**	98	96 %	0.952
555 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	92 %	75266
121 C030	METHYLENE CHLORIDE *		84	34 UG/L	0.680
337 C165	BENZENE		78	8 UG/L	0.873
586 C250	TOTAL XYLEMES		106	11 UG/L	1.216
					1897

* BLANK CORRECTED ON DATA SHEET

RIC
01/13/87 11:43:00
SAMPLE: SCK-MW10-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

511400
DATA: 4U12057C07A #1 SCANS 200 TO 1350
CALI: 4U12057C07A #3

229376.

SPECTRUM DC # -----14

INST ID: 4530

100.0

RIC

200

3:

400

6:40

469

10:00

500

10:20

631

10:40

705

11:00

532

11:20

557

11:40

818

12:00

875

12:20

946

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1031

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1119

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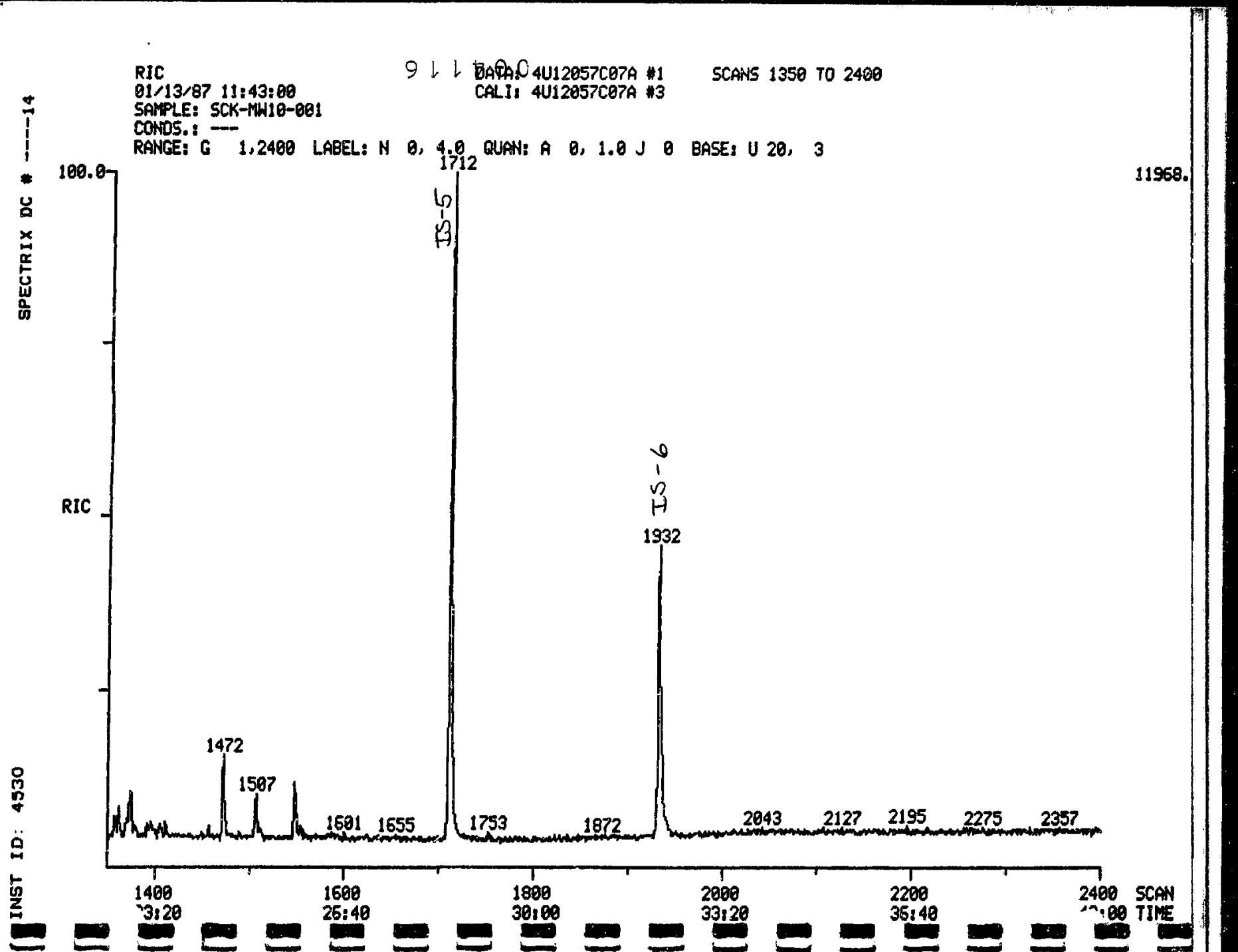
14:40

IS-2

IS-3

IS-4

SCAN
TIME



SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

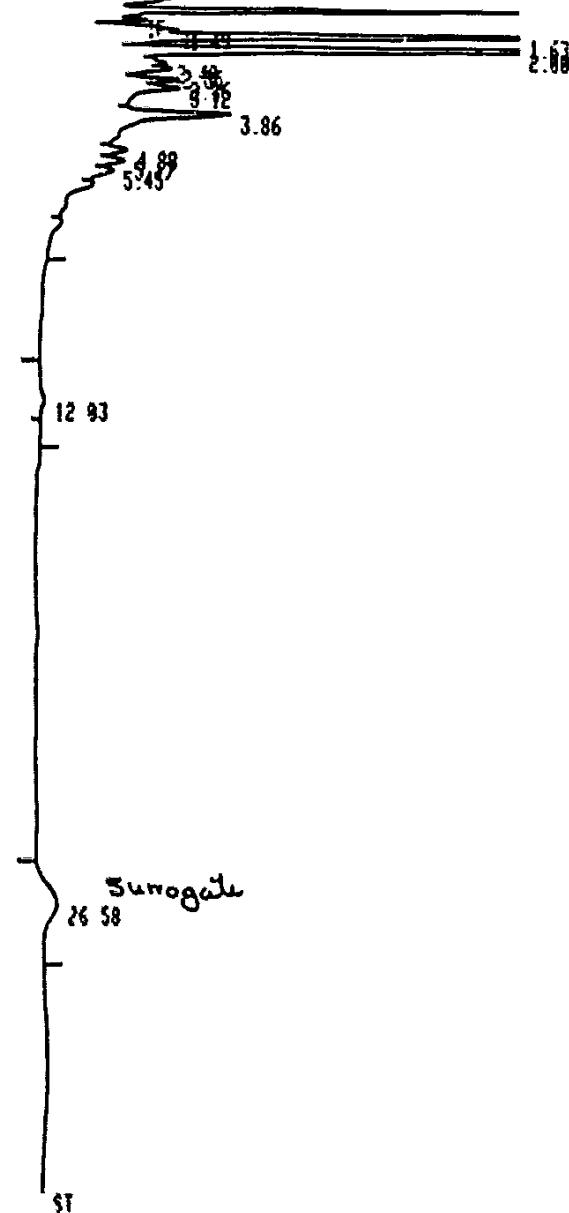
SAMPLE ID.	SCK-MW10-001	INST ID.	4530
FILENAME	4U12057C07A	ANALYST	GU
CLIENT	KEYSTONE	VERIFIED BY	<i>YPT</i>
DATE INJECTED	01/13/87 11:43:00	CORR. FACTOR	20.00
CALIB STD DATE	01/08/87		

SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
532	C130	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	800 UQ/L	1.000	18298
740	C140	NAPHTHALENE-D8 ** IS2 **	136	800 UQ/L	1.000	69080
1031	C150	ACENAPHTHENE-D10 ** IS3 **	164	800 UQ/L	1.000	34881
1272	C160	PHENANTHRENE-D10 ** IS4 **	188	800 UQ/L	1.000	43111
1712	C170	CHRYSENE-D12 ** IS5 **	240	800 UQ/L	1.000	17254
1932	C175	PERYLENE-D12 ** IS6 **	264	800 UQ/L	1.000	10781
743	C450	NAPHTHALENE	128	3413 UQ/L	1.007	29291
858	C470	2-METHYLNAPHTHALENE	142	162 UQ/L	1.159	49331
1036	C550	ACENAPHTHENE	153	683 UQ/L	1.005	36430
1063	C565	DIBENZOFURAN	168	338 UQ/L	1.031	18271
1119	C590	FLUORENE	166	213 UQ/L	1.085	9691
1275	C640	PHENANTHRENE	178	405 UQ/L	1.002	21051
1282	C645	ANTHRACENE	178	50 UQ/L	1.008	2631
1472	C655	FLUORANTHENE	202	56 UQ/L	1.157	2381
507	C715	PYRENE	202	41 UQ/L	0.880	1371

* SURROGATES WERE DILUTED OUT

STAK #: 86-12-057-074

U.A.



DCB _____

SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. SCK-MW10-001
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II.C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 130°C
VOLUME INJ. 1 μl
INSTRUMENT ID. 6000

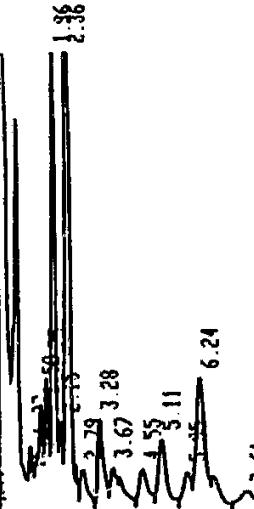
4
00
0

RUN #: 27 JAN/07/87 02:53:31
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%	RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.839	1.49	13980	PY	0.166	1.839
36.939	1.63	280798	D VB	0.094	36.939
46.882	2.00	356380	D PY	0.100	46.882
1.178	2.40	8952	VV	0.187	1.178
1.418	2.55	10778	D VV	0.154	1.418
1.832	2.86	13967	VV	0.156	1.832
1.976	3.12	15021	VV	0.271	1.976
4.890	3.86	31093	VV	0.317	4.890
1.868	4.88	8121	VV	0.231	1.868
1.114	5.12	8467	VV	0.215	1.114
0.883	5.45	6713	VV	0.257	0.883
0.162	12.03	1234	PY	0.641	0.162
0.614	26.58	4666	PB	1.129	0.614

NA.

1001 8C-12-052 - 07A



DCI
SPECTRIX HOUSTON
CASE NO. Key Stone

SAMPLE NO. SEC-MVQ-00
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x 6mm C
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT
COLUMN TEMP. 253°C
COLUMN INJ. 1.0V
INSTRUMENT ID. GC

Surrogate

SI

RUN # 109
WORKFILE ID: C
WORKFILE NAME:

HEIGHT ^Z	HEIGHT ^Z	TYPE	AR/HI	HEIGHT ^Z
1.27	3165	D BY	0.064	1.102
1.58	8928	VY	0.122	3.108
1.75	12682	D VY	0.118	4.415
1.96	5717	D VY	0.105	19.295
2.15	7306	U VY	0.105	2.541
2.36	141288	VY	0.115	49.186
2.79	2614	RP	0.166	0.910
3.28	10931	PY	0.295	3.805
3.67	4614	VY	0.188	1.606
4.55	4491	VY	0.303	1.532
5.11	8499	VY	0.258	2.924
5.85	4297	VY	0.259	1.465
6.24	16550	VY	0.294	5.762
7.61	1595	PB	0.454	0.555
20.62	3402	BY	1.035	1.184

Page 1

Received: 12/12/86

SPECTRIX

REPORT

Work Order # 86-12-007

004120
Results by Sample

SAMPLE ID <u>SCK-MW10-001</u>		SAMPLE # <u>07</u> FRACTIONS: <u>A</u>			
		Date & Time Collected <u>12/12/86</u>		Category	
AG	<10 ug/L	AS	24.7 ug/L	BE	<3 ug/L
CD		CG		CN_SU	<100 ug/L
CR		HG	<0.2 ug/L	NI	<29 ug/L
		PB		SB	<44 ug/L
SE	<5 ug/L	TL	<10 ug/L	ZN	29 ug/L

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-MW13-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

004121

SPECTRIX DC # ---- 8

INST ID 4000

SAMPLE NUMBER: SCK-MW13-601

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205708
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ----
QC REPORT NO.: ----
CONTRACT NO.: ----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V08
DILUTION FACTOR: 1:50

CAS #	COMPOUND	UG/L	Q
C010	CHLOROMETHANE	500 U	0
C015	BROMOMETHANE	500 U	1
C020	VINYL CHLORIDE	500 U	0
C025	CHLOROETHANE	500 U	0
C030	METHYLENE CHLORIDE	250 U	0
C035	ACETONE	500 U	0
C040	CARBON DISULFIDE	250 U	0
C045	1, 1-DICHLOROETHENE	250 U	0
C050	1, 1-DICHLOROETHANE	250 U	0
C055	TRANS-1, 2-DICHLOROETHENE	250 U	0
C060	CHLOROFORM	250 U	0
C065	1, 2-DICHLOROETHANE	250 U	0
C110	2-BUTANONE	500 U	0
C115	1, 1, 1-TRICHLOROETHANE	250 U	0
C120	CARBON TETRACHLORIDE	250 U	0
C125	VINYL ACETATE	500 U	0
C130	BROMODICHLOROMETHANE	250 U	0
C140	1, 2-DICHLOROPROPANE	250 U	0
C145	TRANS-1, 3-DICHLOROPROPENE	250 U	0
C150	TRICHLOROETHENE	250 U	0
C155	DIBROMOCHLOROMETHANE	250 U	0
C160	1, 1, 2-TRICHLOROETHANE	250 U	0
C165	BENZENE	250 U	0
C170	CIS-1, 3-DICHLOROPROPENE	250 U	0
C175	2-CHLOROETHYL VINYL ETHER	500 U	0
C180	BROMOFORM	250 U	0
C210	2-HEXANONE	500 U	0
C205	4-METHYL-2-PENTANONE	500 U	0
C220	TETRACHLOROETHENE	250 U	0
C225	1, 1, 2, 2-TETRACHLOROETHANE	250 U	0
C230	TOLUENE	250 U	0
C235	CHLOROBENZENE	250 U	0
C240	ETHYLBENZENE	250 U	0
C245	STYRENE	250 U	0
C250	TOTAL XYLEMES	250 U	0

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: SCK-MW13-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205708
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86

SEMOVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/18/86.
 DATE ANALYZED: 01/13/87

DATAFILE: 4U12057C08
 DILUTION FACTOR: 200.0....

CAS #	COMPOUND	UG/L	
C315	PHENOL	2000 U	
C325	BIS(2-CHLOROETHYL)ETHER	2000 U	
C330	2-CHLOROPHENOL	2000 U	
C335	1, 3-DICHLOROBENZENE	2000 U	
C340	1, 4-DICHLOROBENZENE	2000 U	
C345	BENZYL ALCOHOL	2000 U	
C350	1, 2-DICHLOROBENZENE	2000 U	
C355	2-METHYLPHENOL	1300 J	
C360	BIS(2-CHLOROISOPROPYL)ETHER	2000 U	
C365	4-METHYLPHENOL	1400 J	
C370	N-NITROSODIISOPROPYLAMINE	2000 U	
C375	HEXACHLOROETHANE	2000 U	
C410	NITROBENZENE	2000 U	
C415	ISOPHORONE	2000 U	
C420	2-NITROPHENOL	2000 U	
C425	2, 4-DIMETHYLPHENOL	6300	
C430	BENZOIC ACID	10000 U	
C435	BIS(2-CHLOROETHOXY)METHANE	2000 U	
C440	2, 4-DICHLOROPHENOL	2000 U	
C445	1, 2, 4-TRICHLOROBENZENE	2000 U	
C450	NAPHTHALENE	26000	
C455	4-CHLOROANILINE	2000 U	
C460	HEXACHLOROBUTADIENE	2000 U	
C465	P-CHLORO-M-CRESOL	2000 U	
C470	2-METHYLNAPHTHALENE	1400 J	
C510	HEXACHLOROCYCLOPENTADIENE	2000 U	
C515	2, 4, 6-TRICHLOROPHENOL	2000 U	
C520	2, 4, 5-TRICHLOROPHENOL	10000 U	
C525	2-CHLORONAPHTHALENE	2000 U	
C530	2-NITROANILINE	10000 U	
C535	DIMETHYL PHTHALATE	2000 U	
C540	ACENAPHTHYLENE	2000 U	
C545	3-NITROANILINE	10000 U	
C550	ACENAPHTHENE	940 J	
C555	2, 4-DINITROPHENOL	10000 U	
C560	4-NITROPHENOL	10000 U	
C565	DIBENZOFURAN	510 J	
C570	2, 4-DINITROTOLUENE	2000 U	
C575	2, 6-DINITROTOLUENE	2000 U	

004123

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-MW13-001

MIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C08

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	2000 U
C585	4-CHLOROPHENYL PHENYL ETHER	2000 U
C590	FLUORENE	390 U
C595	4-NITROANILINE	10000 U
C610	4,6-DINITRO-2-METHYLPHENOL	10000 U
C615	N-NITROSODIPHENYLAMINE	2000 U
C625	4-BROMOPHENYL PHENYL ETHER	2000 U
C630	HEXACHLOROBENZENE	2000 U
C635	PENTACHLOROPHENOL	10000 U
C640	PHENANTHRENE	380 U
C645	ANTHRACENE	2000 U
C650	DI-N-BUTYL PHTHALATE	2000 U
C655	FLUORANTHENE	2000 U
C715	PYRENE	2000 U
C720	BUTYL BENZYL PHTHALATE	2000 U
C725	3,3'-DICHLOROBENZIDINE	4000 U
C730	BENZO(A)ANTHRACENE	2000 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	2000 U
C740	CHRYSENE	2000 U
C760	DI-N-OCTYL PHTHALATE	2000 U
C765	BENZO(B)FLUORANTHENE	2000 U
C770	BENZO(K)FLUORANTHENE	2000 U
C775	BENZO(A)PYRENE	2000 U
C780	INDENO(1,2,3-CD)PYRENE	2000 U
C785	DIBENZO(A,H)ANTHRACENE	2000 U
C790	BENZO(GHI)PERYLENE	2000 U

004124

U = UNDETECTED AT THE LISTED DETECTION LIMIT

U = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-08A

DC No: KEYSTONE
SMO No: SCK-MW13-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/18/86
Date Analyzed: 1/7/87
Dilution Factor: 1
Percent Moisture(decantated):

GPC Cleanup Yes No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOR-1016	0.50 U
11104-28-2	AROCLOR-1221	0.50 U
11141-16-5	AROCLOR-1232	0.50 U
53469-21-9	AROCLOR-1242	0.50 U
12672-29-6	AROCLOR-1248	0.50 U
11097-69-1	AROCLOR-1254	1.0 U
11096-82-5	AROCLOR-1260	1.0 U

004125

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-MW13-00

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO. : -----

ANALYST: NOT

DATAFILE: 4U12057V08

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
			6	0
			2	UG/
			1	
			4	
	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.		0	0

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DG # ---- 8

SAMPLE: SCK-MW13

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12057C08

B. TENTATIVELY IDENTIFIED COMPOUNDS

IAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
95-13-6	1H-INDENE	568	937	1400 J
498-81-7	CYCLOHEXANEMETHANOL, . ALPHA., . ALPHA., 4-TR	696	858	1100 J
526-75-0	PHENOL, 2,3-DIMETHYL-	726	908	3900 J
95-15-8	BENZO[C,B]THIOPHENE	749	844	2800 J
99-89-8	PHENOL, 4-(1-METHYLETHYL)-	789	877	1100 J
1515-95-3	BENZENE, 1-ETHYL-4-METHOXY-	800	876	900 J
3855-26-3	PHENOL, 2-ETHYL-4-METHYL-	822	849	1800 J
2471-83-2	1H-INDENE, 1-ETHYLLIDENE-	874	793	1100 J
6351-10-6	1H-INDEN-1-OL, 2,3-DIHYDRO-	880	755	880 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	946		960 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1101		2500 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 46009

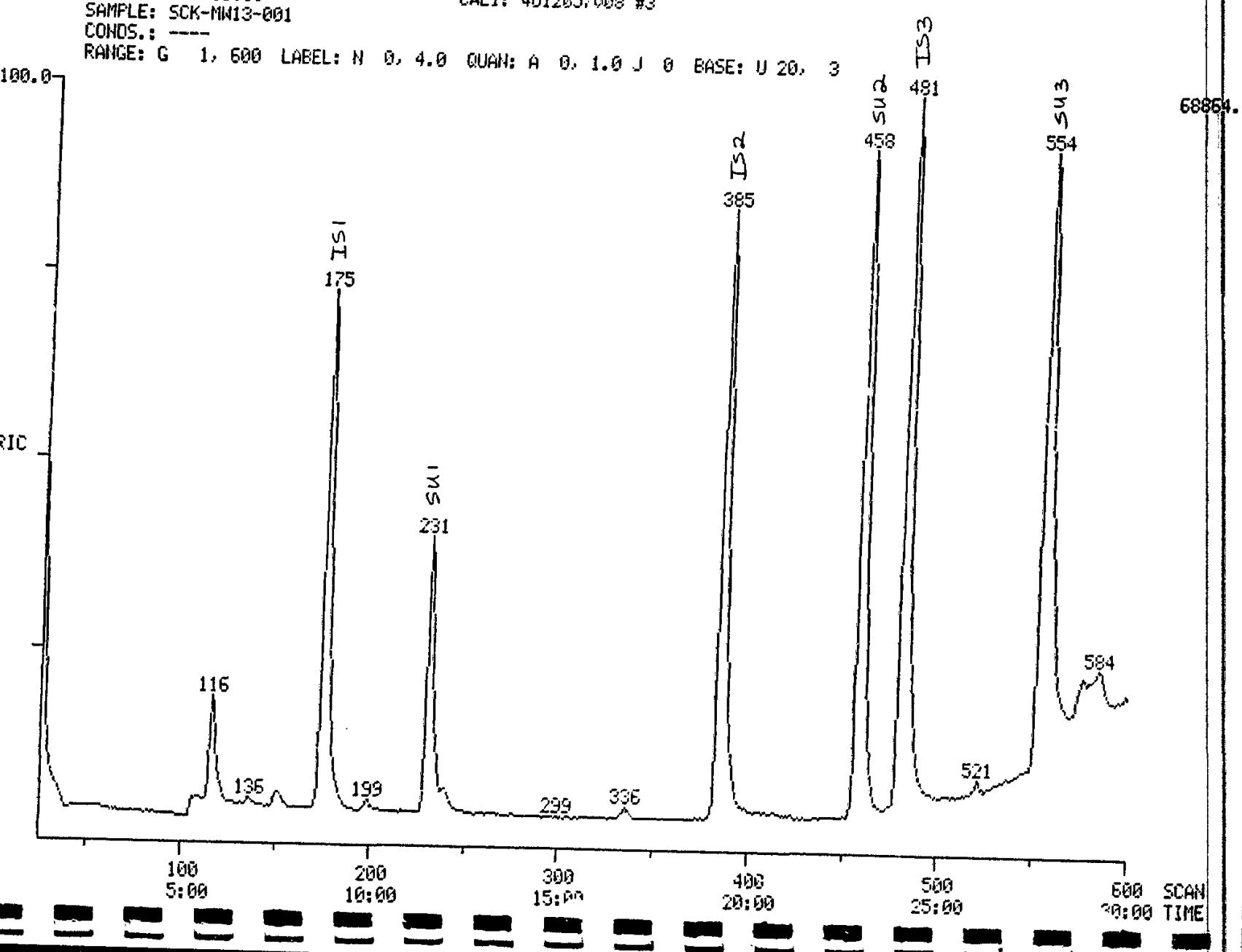
SPECTRUM DC # 14

RIC
12/19/86 15:59:00
SAMPLE: SCK-MN13-001
COND.: -----

DATA: 4U12057V08 #1
CALI: 4U12057V08 #3

SCANS 25 TO 600

RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3



SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID. SCK-MW13-001

FILENAME 4U12057V08 INST 4000

CLIENT KEYSTONE ANALYST NOT

DATE INJECTED 12/19/86 15:59:00 VERIFIED BY *Not*

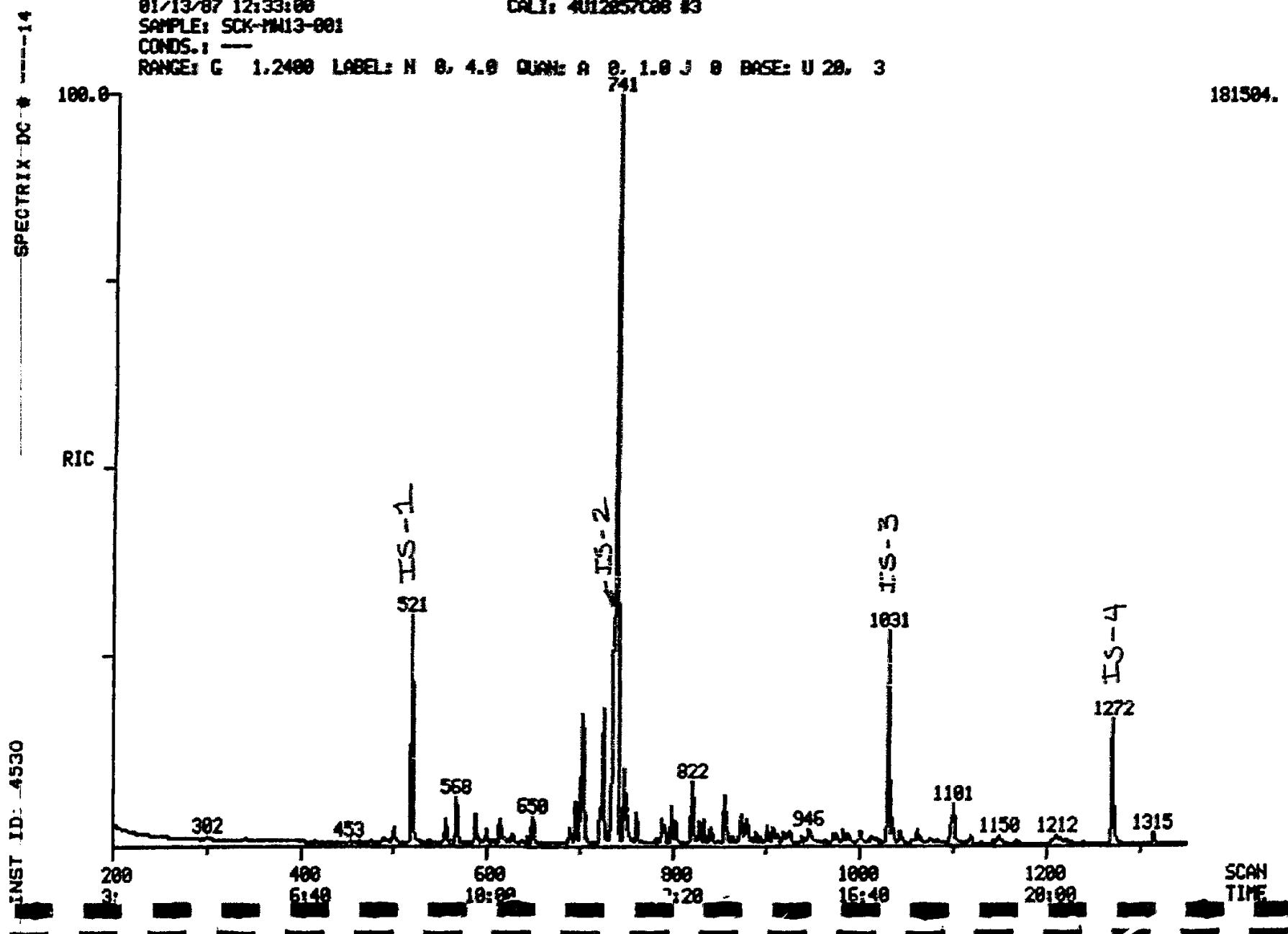
CALIB STD DATE 12-14-86 CORR. FACTOR 50.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
175 CI01	BROMOCHLOROMETHANE	**IS1**	128	2500 UG/L	1.000 21001
385 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	2500 UG/L	1.000 97316
481 CI20	D5-CHLOROBENZENE	**IS3**	117	2500 UG/L	1.000 87826
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	86 %	1.320 27373
458 CS05	D8-TOLUENE	**SU2**	98	93 %	0.952 83949
554 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	86 %	1.152 55190
115 C030	METHYLENE CHLORIDE *		84	395 UG/L	0.663 61610

* BLANK CORRECTED ON DATA SHEET

RIC
01/13/87 12:33:00
SAMPLE: SCK-MN13-001
COND.S.:
RANGE: G 1.2400 LABEL: N 8, 4.9 QUAN: A 8, 1.0 J 8 BASE: U 20, 3

0 < 1 μ g C
DATE: 4U12857088 #1
CALI: 4U12857088 #3
SCANS 200 TO 1350

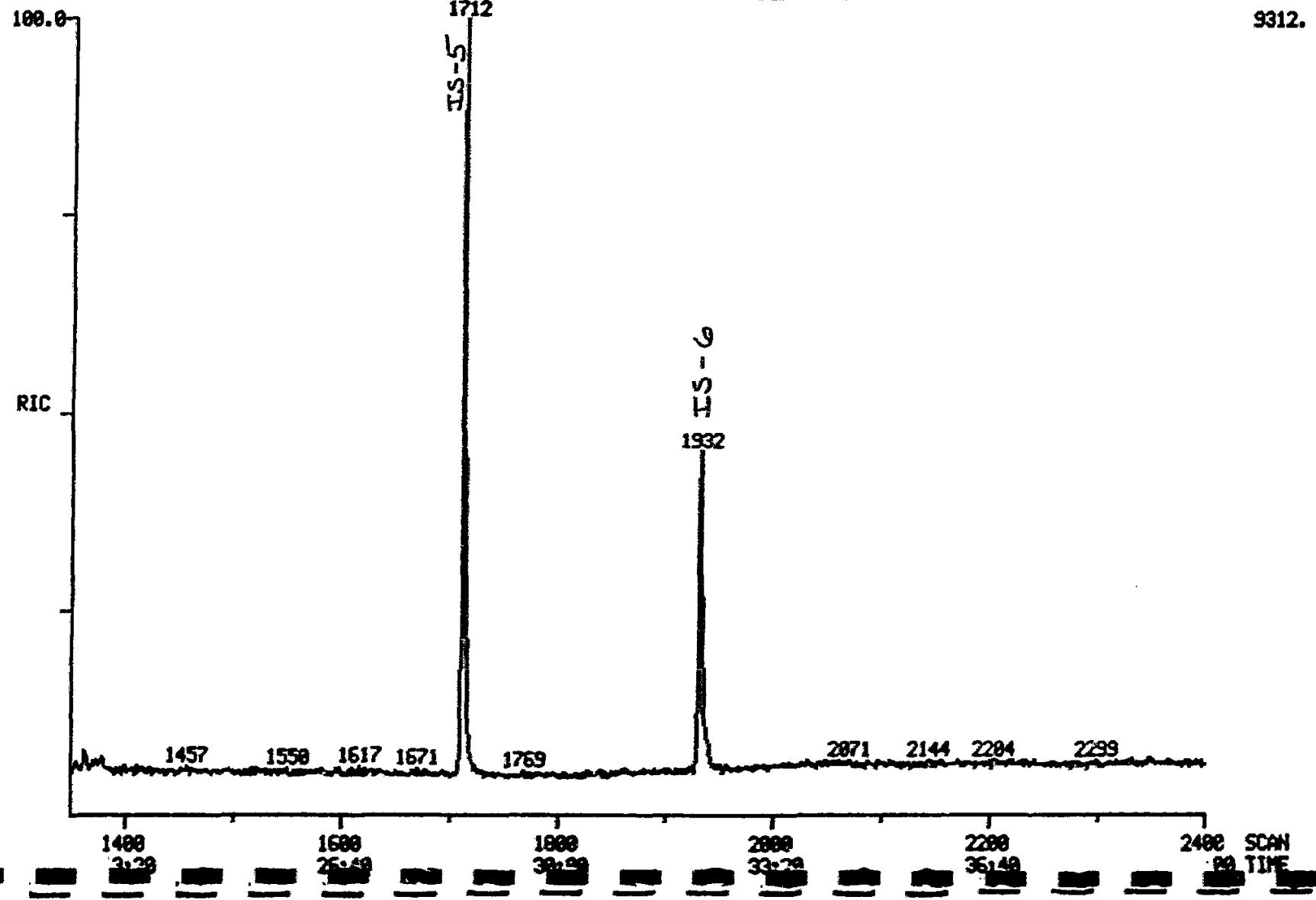


RIC
01/13/87 12:33:00
SAMPLE: SCK-MN13-901
COND.: —

L C L + 0.0
DATA: 4U12657088 #1
CALI: 4U12657088 #3
SCANS 1358 TO 2490

RANGE: G 1.2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

SPECTRUM DC # -14



004131

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	SCK-MW13-001		
FILENAME	4U12057C08	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/13/87 12:33:00	VERIFIED BY	<i>JPM</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	200.00

SCANS		SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
521	C130	1,4-DICHLOROBENZEND-D4 ** IS1 **	192	8000 UG/L	1.000	1977.
737	C140	NAPHTHALENE-D8 ** IS2 **	136	8000 UG/L	1.000	7132
1031	C150	ACENAPHTHENE-D10 ** IS3 **	164	8000 UG/L	1.000	3376
1272	C160	PHENANTHRENE-D10 ** IS4 **	198	8000 UG/L	1.000	3587
1712	C170	CHRYSENE-D12 ** IS5 **	240	8000 UG/L	1.000	1468
1932	C175	PERYLENE-D12 ** IS6 **	264	8000 UG/L	1.000	903
589	C355	2-METHYLPHENOL	108	1315 UG/L	1.129	481
613	C365	4-METHYLPHENOL	108	1435 UG/L	1.180	518
703	C425	2,4-DIMETHYLPHENOL	122	6258 UG/L	0.954	2015
741	C450	NAPHTHALENE	128	26013 UG/L	1.005	3054
857	C470	2-METHYLNAPHTHALENE	142	1372 UG/L	1.163	816
1035	C550	ACENAPHTHENE	153	943 UG/L	1.004	353
1062	C565	DIBENZOFURAN	168	510 UG/L	1.030	266
1119	C590	FLUORENE	166	391 UG/L	1.085	172
275	C640	PHENANTHRENE	178	380 UG/L	1.002	164

* SURROGATES WERE DILUTED OUT

INST ID: 4530

SPECTRIX DC # ---19

1145
SAMPLE

ANALYST: GU

C10.H12.0
M WT 1145
B PK 133
RANK 1
7166
PUR 699

C11.H16
M WT 1145
B PK 133
RANK 1
7165
PUR 688

C18.H12.0
M WT 1145
B PK 133
RANK 3
7154
PUR 675

M/Z

LIBRARY SEARCH 0041400
01/13/87 12:33:00 + 15:46
SAMPLE: SCK-MN13-001
COND.:
ENHANCED (S 158 2N 0T)

DATA: 4U12057C08 # 945
CALL: 4U12057C08 # 3
BASE M/Z: 133
RIC: 2371.

004133

SPECTRUM DC # 19

1309
SAMPLE

ANALYST: GU

C9.H9.O.N
M WT 147
B PK 118
RANK 1
7103
PUR 663

1309
M WT 147
B PK 147
RANK 2
7082
PUR 653

1309
M WT 147
B PK 147
RANK 3
7080
PUR 645

INST ID: 4530

M/Z

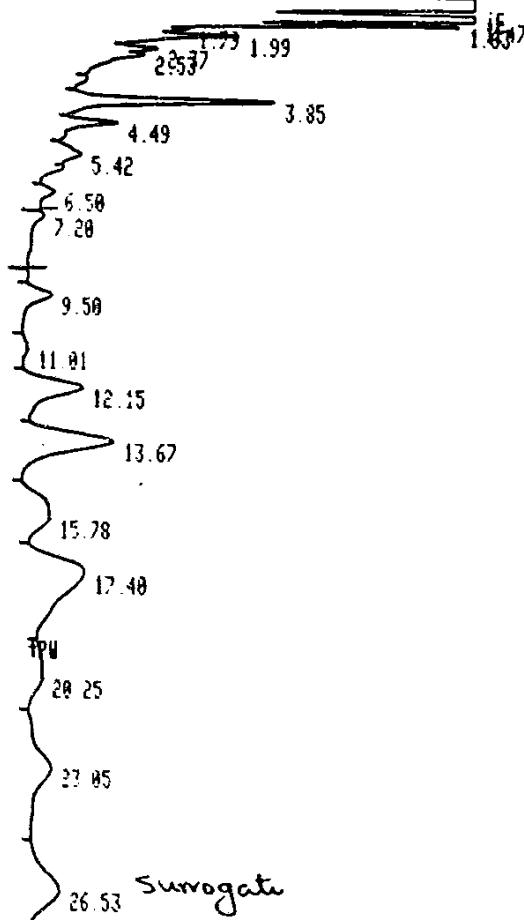
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LIBRARY SEARCH
01/13/87 12:33:00 + 18:21
SAMPLE: SCK-MN13-001
COND.: --
ENHANCED (S 15B 2N 0T)

004134
DATA: 4U12057C08 #1101
CALI: 4U12057C08 # 3
BASE M/Z: 147
RIC: 7727.

86-12-057-08A

NA.



RUN # 28 JAN/07/87 03:29:48
WORKFILE ID: C
WORKFILE NAME:

HEIGHTS

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.47	133222	BY	0.153	31.420
1.53	66244	D VP	0.084	15.623
1.79	3950	D PV	0.089	0.932
1.99	22807	D VF	0.132	5.379
2.37	9938	PV	0.137	2.344
2.53	8511	D VP	0.181	2.002
3.85	52838	PV	0.280	12.462
4.49	14525	VV	0.260	3.426
5.42	8345	VV	0.414	1.968
6.58	3741	PB	0.272	0.882
9.50	7080	VP	0.415	1.670
11.01	1521	PV	0.492	0.359
12.15	16184	VV	0.619	3.817
13.67	24196	VV	0.642	5.685
15.78	7577	VV	1.142	1.787
17.40	17097	VV	1.571	4.832
20.25	6032	VV	1.665	1.423
23.05	8312	VV	1.950	1.960
26.53	10703	VP	1.846	2.524
33.96	1271	I PP	1.333	0.306

DC#
SPECTRIX HOUSTON

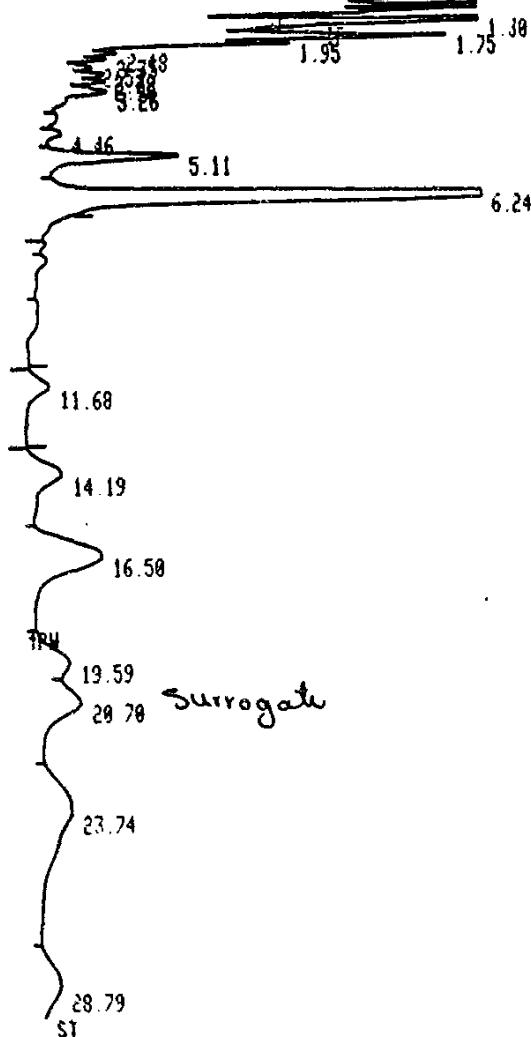
CASE NO. Keystone

SAMPLE NO. SCK-MW13-0
PESTICIDE FRACTION
2ND COLUMN CONFIRMA

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 184°C
VOLUME INJ. 1 uL
INSTRUMENT ID. 6800

04

STK # 86-12-057-08A



DC#
SPECTRIX HOUSTON

CASE NO. Key Stone

SAMPLE NO. SKC-MW13-6
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x 6mm C
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 70.3°C
VOLUME INJ. 1 μl
INSTRUMENT ID. 5600
O
O

RUN # 110 JAN/12/87 18:21:26
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%	RT	HEIGHT	TYPE	AR/HT	HEIGHT%
53.110	253576	0	BB	0.084	1.38
6.622	31615	8	BV	0.139	1.75
3.873	18490	0	VP	0.104	1.95
0.554	2646	0	PV	0.105	2.18
0.584	2789	0	VP	0.091	2.35
0.488	2291	PV		0.132	2.57
0.892	4260	VV		0.156	2.78
0.979	4673	0	VV	0.154	2.96
1.188	5636	V	VP	0.274	3.26
0.328	1568	P	PP	0.210	4.46
3.452	16479	P	PP	0.244	5.11
20.799	99304	P	PB	0.263	6.24
0.563	2686	B	BB	0.507	11.68
0.901	4300	B	BV	0.919	14.19
1.989	9114	V	VV	1.091	16.50
0.989	4724	V	VV	0.983	19.59
1.270	6065	V	VV	1.231	20.70
0.962	4592	V	VV	2.457	23.74
0.554	2643	I	VP	1.361	28.79

Page 1

Received: 12/12/86

SPECTRIX

REPORT

Work Order # 86-12-03/

004137 Results by Sample

SAMPLE ID SCK-MW13-001		SAMPLE # 08 FRACTIONS: A									
		Date & Time Collected 12/12/86 Category									
AG	<10	AS	97.9	BE	<3	CD	B	CN_SU	<100	CR	10
CU	<10	FE	5110	HG	<0.2	NI	<29	PB	<5	SB	<44
SE	<5	TL	<10	ZN	29						

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-MW22-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

004138

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: SCK-MW22-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861205709
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *YPH*

CASE NO. : -----
QC REPORT NO. : -----
CONTRACT NO. : -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/19/86
DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V09

DILUTION FACTOR: 1:1.....

CAS #	COMPOUND	UG/L	Q	M
C010	CHLOROMETHANE	10 U		1
C015	BROMOMETHANE	10 U		4
C020	VINYL CHLORIDE	10 U		0
C025	CHLOROETHANE	10 U		0
C030	METHYLENE CHLORIDE	40		
C035	ACETONE	10 U		
C040	CARBON DISULFIDE	5 U		
C045	1, 1-DICHLOROETHENE	5 U		
C050	1, 1-DICHLOROETHANE	5 U		
C055	TRANS-1, 2-DICHLOROETHENE	5 U		
C060	CHLOROFORM	5 U		
C065	1, 2-DICHLOROETHANE	5 U		
C110	2-BUTANONE	10 U		
C115	1, 1, 1-TRICHLOROETHANE	5 U		
C120	CARBON TETRACHLORIDE	5 U		
C125	VINYL ACETATE	10 U		
C130	BROMODICHLOROMETHANE	5 U		
C140	1, 2-DICHLOROPROPANE	5 U		
C145	TRANS-1, 3-DICHLOROPROPENE	5 U		
C150	TRICHLOROETHENE	5 U		
C155	DIBROMOCHLOROMETHANE	5 U		
C160	1, 1, 2-TRICHLOROETHANE	5 U		
C165	BENZENE	5 U		
C170	CIS-1, 3-DICHLOROPROPENE	5 U		
C175	2-CHLOROETHYL VINYLETHER	10 U		
C180	BROMOFORM	5 U		
C210	2-HEXANONE	10 U		
C205	4-METHYL-2-PENTANONE	10 U		
C220	TETRACHLOROETHENE	5 U		
C225	1, 1, 2, 2-TETRACHLOROETHANE	5 U		
C230	TOLUENE	5 U		
C235	CHLOROBENZENE	5 U		
C240	ETHYLBENZENE	5 U		
C245	STYRENE	5 U		
C250	TOTAL XYLENES	5 U		

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX UC # ---- 8
SAMPLE NUMBER: SCK-MW22-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205709
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86.

SEMIVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: ... (2/10/86)
 DATE ANALYZED: 01/13/87

DATAFILE: 4U12057C09
 DILUTION FACTOR: 1.0.....

CAS #	COMPOUND	UG/L	
C315	PHENOL	10 U	
C325	BIS(2-CHLOROETHYL)ETHER	10 U	
C330	2-CHLOROPHENOL	10 U	
C335	1, 3-DICHLOROBENZENE	10 U	
C340	1, 4-DICHLOROBENZENE	10 U	
C345	BENZYL ALCOHOL	10 U	
C350	1, 2-DICHLOROBENZENE	10 U	
C355	2-METHYLPHENOL	10 U	
C360	BIS(2-CHLOROISOPROPYL)ETHER	10 U	
C365	4-METHYLPHENOL	10 U	
C370	N-NITROSODIPropylamine	10 U	
C375	HEXACHLOROETHANE	10 U	
C410	NITROBENZENE	10 U	
C415	ISOPHORONE	10 U	
C420	2-NITROPHENOL	10 U	
C425	2, 4-DIMETHYLPHENOL	10 U	
C430	BENZOIC ACID	50 U	
C435	BIS(2-CHLOROETHOXY)METHANE	10 U	
C440	2, 4-DICHLOROPHENOL	10 U	
C445	1, 2, 4-TRICHLOROBENZENE	10 U	
C450	NAPHTHALENE	10 U	
C455	4-CHLOROANILINE	10 U	
C460	HEXACHLOROBUTADIENE	10 U	
C465	P-CHLORO-M-CRESOL	10 U	
C470	2-METHYLNAPHTHALENE	10 U	
C510	HEXACHLOROCYCLOPENTADIENE	10 U	
C515	2, 4, 6-TRICHLOROPHENOL	10 U	
C520	2, 4, 5-TRICHLOROPHENOL	50 U	
C525	2-CHLORONAPHTHALENE	10 U	
C530	2-NITROANILINE	50 U	
C535	DIMETHYL PHTHALATE	10 U	
C540	ACENAPHTHYLENE	10 U	
C545	3-NITROANILINE	50 U	
C550	ACENAPHTHENE	10 U	
C555	2, 4-DINITROPHENOL	50 U	
C560	4-NITROPHENOL	50 U	
C565	DIBENZOFURAN	10 U	
C570	2, 4-DINITROTOLUENE	10 U	
C575	2, 6-DINITROTOLUENE	10 U	

004140

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-MN22-001

IN VOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12057C09

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	10 U
C585	4-CHLOROPHENYL PHENYL ETHER	10 U
C590	FLUORENE	10 U
C595	4-NITROANILINE	50 U
C610	4,6-DINITRO-2-METHYLPHENOL	50 U
C615	N-NITROSODIPHENYLAMINE	10 U
C625	4-BROMOPHENYL PHENYL ETHER	10 U
C630	HEXACHLOROBENZENE	10 U
C635	PENTACHLOROPHENOL	50 U
C640	PHENANTHRENE	10 U
C645	ANTHRACENE	10 U
C650	DI-N-BUTYL PHTHALATE	10 U
C655	FLUORANTHENE	10 U
C715	PYRENE	10 U
C720	BUTYL BENZYL PHTHALATE	10 U
C725	3,3'-DICHLOROBENZIDINE	20 U
C730	BENZO(A)ANTHRACENE	10 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	10 U
C740	CHRYSENE	10 U
C760	DI-N-OCTYL PHTHALATE	10 U
C765	BENZO(B)FLUORANTHENE	10 U
C770	BENZO(K)FLUORANTHENE	10 U
C775	BENZO(A)PYRENE	10 U
C780	INDENO(1,2,3-CD)PYRENE	10 U
C785	DIBENZO(A,H)ANTHRACENE	10 U
C790	BENZO(GHI)PERYLENE	10 U

004141

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-09A

DC No: KEYSTONE
SMO No: SCK-MW22-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low GPC Cleanup Yes ✓ No
Date Extracted: 12/18/86 Sep. Funnel Ext. Yes
Date Analyzed: 1/7/87 Cont. Liquid Ext. Yes

Dilution Factor: 1

Percent Moisture(decant):

CAS #	Compound	ug/L	
319-84-6	ALPHA-BHC	0.05	U
319-85-7	BETA-BHC	0.05	U
319-86-8	DELTA-BHC	0.05	U
58-89-9	GAMMA-BHC (LINDANE)	0.05	U
76-44-8	HEPTACHLOR	0.05	U
309-00-2	ALDRIN	0.05	U
1024-57-3	HEPTACHLOR EPOXIDE	0.05	U
959-98-8	ENDOSULFAN I	0.05	U
60-57-1	DIELDRIN	0.10	U
72-55-9	4,4'-DDE	0.10	U
72-20-8	ENDRIN	0.10	U
33213-65-9	ENDOSULFAN II	0.10	U
72-54-8	4,4'-DDD	0.10	U
7421-93-4	ENDRIN ALDEHYDE	0.10	U
1031-07-8	ENDOSULFAN SULFATE	0.10	U
50-29-3	4,4'-DDT	0.10	U
72-43-5	METHOXYPHOR	0.50	U
53494-70-5	ENDRIN KETONE	0.10	U
57-74-9	CHLORDANE	0.50	U
8001-35-2	TOXAPHENE	1.0	U
12674-11-2	AROCLOTR-1016	0.50	U
11104-28-2	AROCLOTR-1221	0.50	U
11141-16-5	AROCLOTR-1232	0.50	U
53469-21-9	AROCLOTR-1242	0.50	U
12672-29-6	AROCLOTR-1248	0.50	U
11097-69-1	AROCLOTR-1254	1.0	U
11096-82-5	AROCLOTR-1260	1.0	U

Vs = 500 = Volume of water extracted (ml)

Vt = 5000 = Volume of total extract (ul)

Vi = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-MW22-000

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO. : -----

ANALYST: CKT

DATAFILE: 4U12057V09

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
			M	
			4	UG/
			1	
	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.		0	
			0	
			4	
			0	
			0	

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4330

SPECTRIX DC # --- 8

SAMPLE: SCK-MW22

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: QU

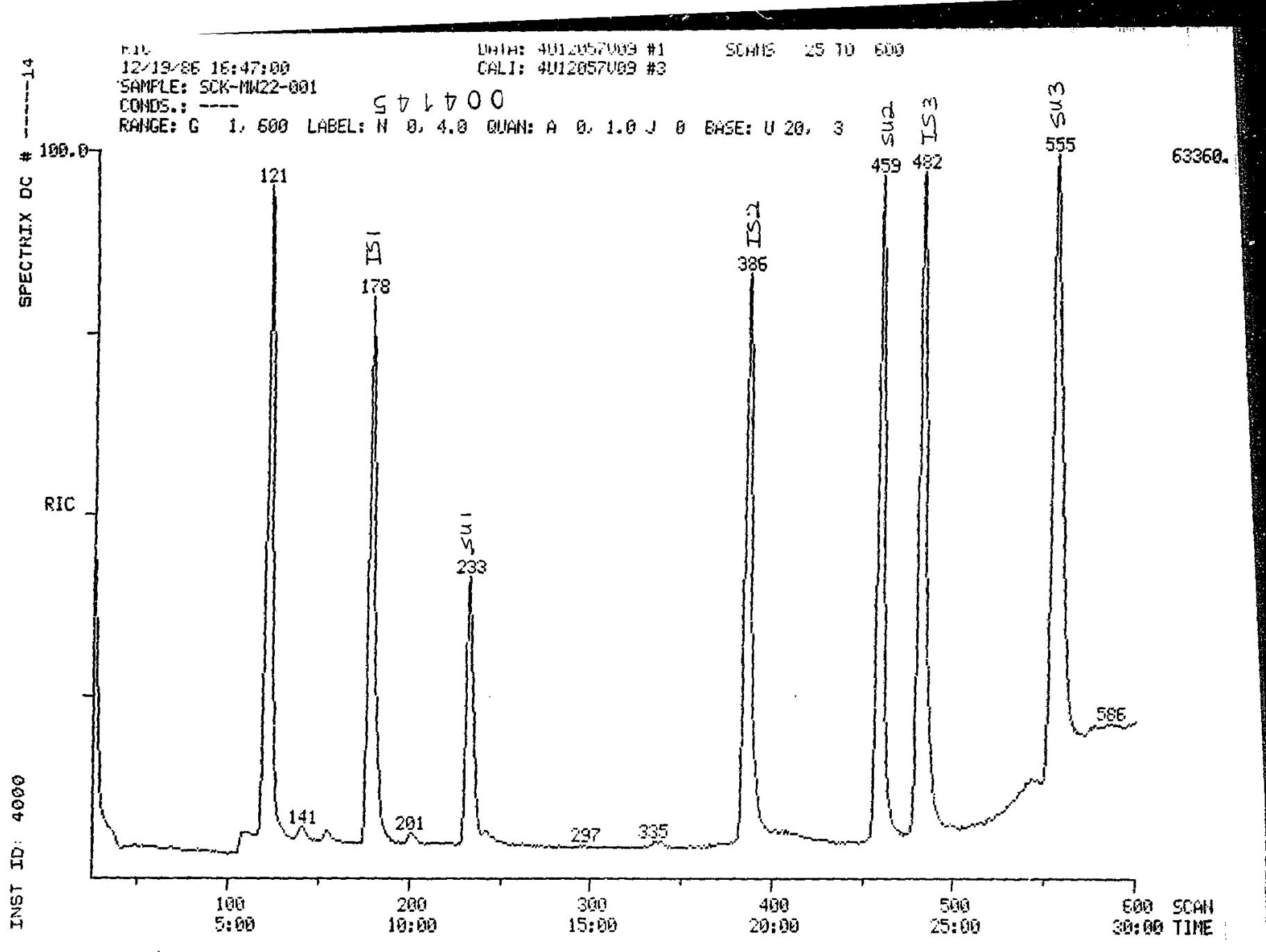
DATAFILE: 4U12057C09

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
110-12-3	2-HEXANONE, 5-METHYL-	* 289	936	44 13 J 41 400

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK



INST ID: 4530

SPECTRIX DC # ----14

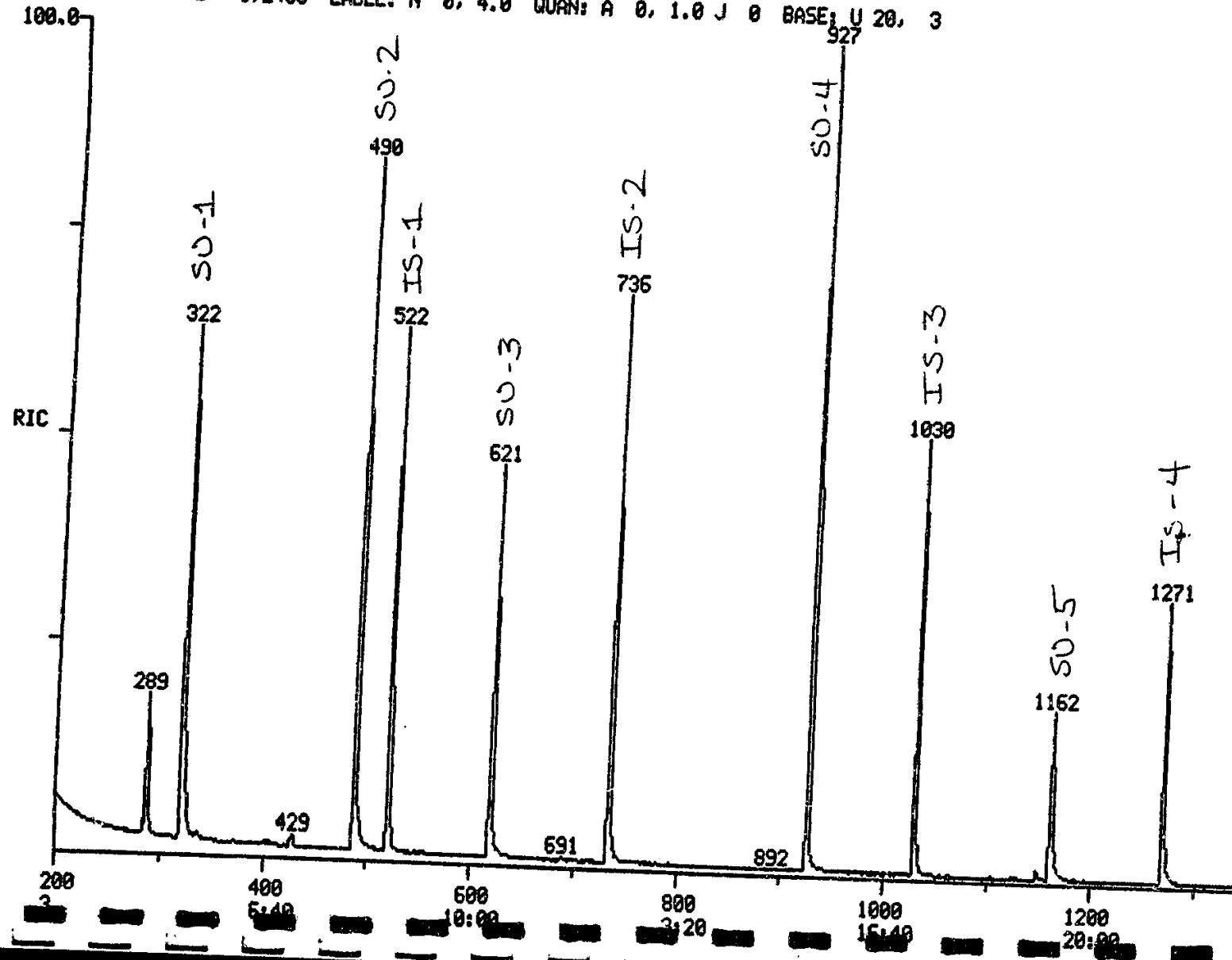
100.0

RIC
01/13/87 13:24:00
SAMPLE: SCK-MW22-001
COND.: --
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

004146

DATA: 4U12057C09 #1
CALI: 4U12057C09 #3

SCANS 200 TO 1350



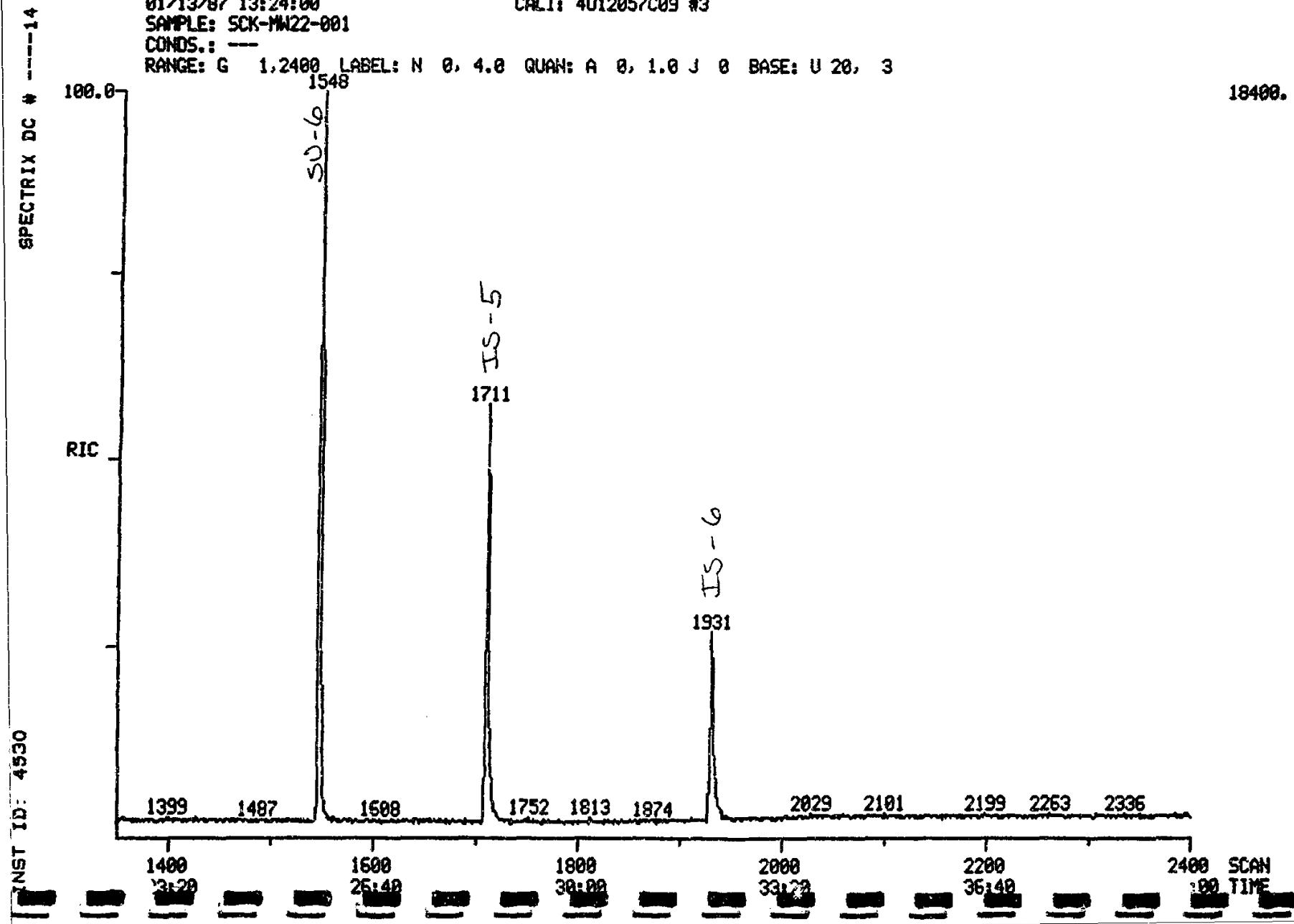
004146

78000.

SCAN
TIME

RIC
01/13/87 13:24:00
SAMPLE: SCK-MN22-001
CONDNS: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

0004147
DATA: 4U12057C09 #1
CALI: 4U12057C09 #3
SCANS 1350 TO 2400



SPECTRIX DC # ----14

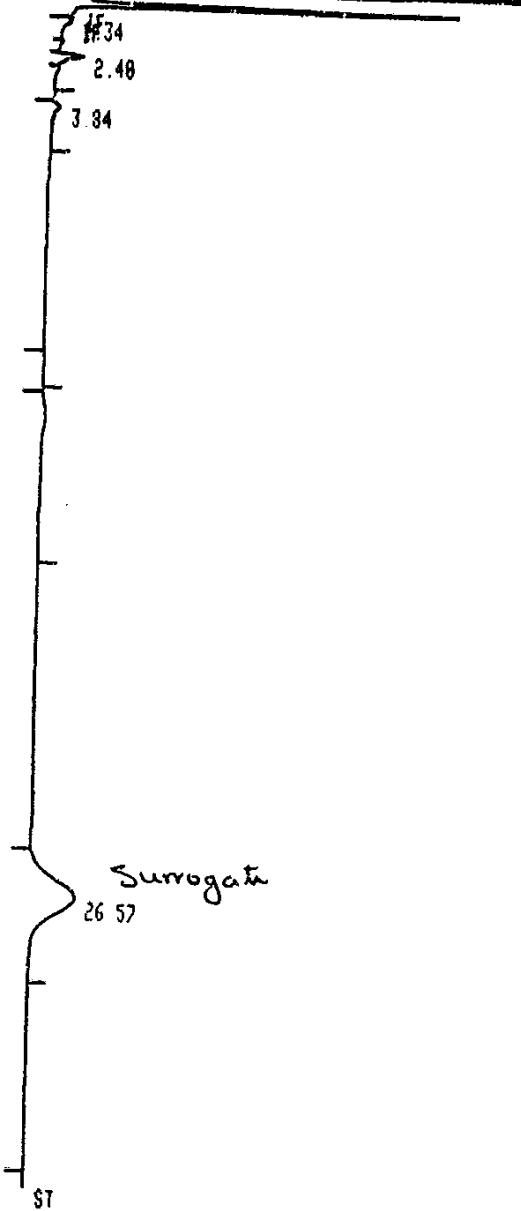
SEMIVOLATILE WATER REPORT

SAMPLE ID.	SCK-MW22-001		
ILENAME	4U12057C09	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/13/87 13:24:00	VERIFIED BY	ypm
CALIB STD DATE	01/08/87	CORR. FACTOR	1.00

SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
322	CI30	1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	40 UG/L	1.000	1606:
736	CI40	NAPHTHALENE-D8 ** IS2 **	136	40 UG/L	1.000	5799:
030	CI50	ACENAPHTHENE-D10 ** IS3 **	164	40 UG/L	1.000	2969:
1271	CI60	PHENANTHRENE-D10 ** IS4 **	188	40 UG/L	1.000	03421:
711	CI70	CHRYSENE-D12 ** IS5 **	240	40 UG/L	1.000	4624:
931	CI75	PERYLENE-D12 ** IS6 **	264	40 UG/L	1.000	1074:
322	CS50	2-FLUOROPHENOL ** SU1 **	112	73 %	0.617	3907:
490	CS43	PHENOL-D5 ** SU2 **	99	75 %	0.939	5928:
162	CS55	2, 4, 6-TRIBROMOPHENOL ** SU5 **	330	67 %	1.128	575:
621	CS20	NITROBENZENE-D5 ** SU3 **	82	96 %	0.844	Q219:
927	CS25	2-FLUOROBIPHENYL ** SU4 **	172	96 %	0.900	7375:
547	CS30	TERPHENYL-D14 ** SU6 **	244	115 %	0.904	2468:

STK# 86-12-057-09A

NA.



ST

DC#
SPECTRIX HOUSTON
CASE NO. Keystone

SAMPLE NO. SCK-MW22-0
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 137°C
VOLUME INJ. 1 AL
INSTRUMENT ID. 6000

0
0

RUN # 29
WORKFILE ID: C JAN/97/87 04:06:05
WORKFILE NAME:

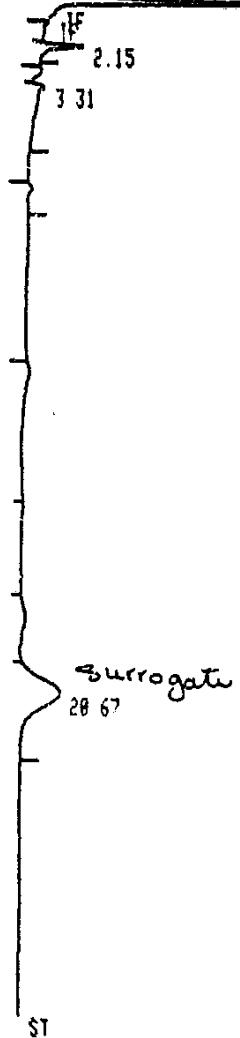
HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.34	1528	D PP	0.111	2.146
2.49	6498	PV	0.147	30.389
3.84	1740	BB	0.210	8.137
26.57	11617	BB	1.208	54.328

TOTAL HGT= 21383
FUL FACTOR= 1.0000E+00

RT# 12-057-09A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. SCK-MWZ-0
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 50°C
VOLUME INJ. 1µl
INSTRUMENT ID. 600C

O

RUN # 111 JAN/12/87 18:52:43
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
2.15	5533	PB	0.142	45.872
3.31	1596	PB	0.615	13.881
28.67	5147	PB	0.964	41.927

TOTAL HIGHT= 12276
MUL FACTOR= 1.0000E+00

Page 1
Received: 12/12/86

SPECTRIX

REPORT

Work Order # 86-12-057

1510 Results by Sample

SAMPLE ID <u>SCK-MW22-001</u>		SAMPLE # 09 FRACTIONS: A									
		Date & Time Collected <u>12/12/86</u>								Category _____	
AG	<10	AS	<10	BE	<3	CD	<5	CN SU	<100	CR	<9
CU	<10	FE	<30	HG	<0.2	NI	<29	PB	<5	SB	<44
SE	<5	TL	<10	ZN	18						

CHECKLIST

SECTION II - Sample Data Packets

- Sample No: TRIP BLANK
- Organic Analysis Data Sheets (OADS)
 - VOA
 - SV
 - NA Pest.
 - Tentatively Identified Compounds
 - VOA
 - SV
 - Raw Data
 - VOA chromatogram
 - VOA quantitation report (surrogates)
 - NA SV chromatogram
 - NA SV quantitation report (surrogates)
 - NA Pest. chromatogram and integrator print out
 - Inorganic Analysis Data Sheets (IADS)

004152

INST ID 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: TRIP BLANK

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861205710
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *YPH*

CASE NO.: -----
 QC REPORT NO.: -----
 CONTRACT NO.: -----
 DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/19/86
 DATE ANALYZED: 12/19/86

DATAFILE: 4U12057V10

DILUTION FACTOR: 1:1

CAS #	COMPOUND	UG/L	LOD
C010	CHLOROMETHANE	10	U
C015	BROMOMETHANE	10	U
C020	VINYL CHLORIDE	10	U
C025	CHLOROETHANE	10	U
C030	METHYLENE CHLORIDE	10	U
C035	ACETONE	44	
C040	CARBON DISULFIDE	10	U
C045	1, 1-DICHLOROETHENE	5	U
C050	1, 1-DICHLOROETHANE	5	U
C055	TRANS-1, 2-DICHLOROETHENE	5	U
C060	CHLOROFORM	5	U
C065	1, 2-DICHLOROETHANE	5	U
C110	2-BUTANONE	5	U
C115	1, 1, 1-TRICHLOROETHANE	10	U
C120	CARBON TETRACHLORIDE	5	U
C125	VINYL ACETATE	5	U
C130	BROMODICHLOROMETHANE	5	U
C140	1, 2-DICHLOROPROPANE	10	U
C145	TRANS-1, 3-DICHLOROPROPENE	5	U
C150	TRICHLOROETHENE	5	U
C155	DIBROMOCHLOROMETHANE	5	U
C160	1, 1, 2-TRICHLOROETHANE	5	U
C165	BENZENE	5	U
C170	CIS-1, 3-DICHLOROPROPENE	5	U
C175	2-CHLOROETHYL VINYL ETHER	5	U
C180	BROMOFORM	10	U
C210	2-HEXANONE	5	U
C205	4-METHYL-2-PENTANONE	10	U
C220	TETRACHLOROETHENE	10	U
C225	1, 1, 2, 2-TETRACHLOROETHANE	5	U
C230	TOLUENE	5	U
C235	CHLOROBENZENE	5	U
C240	ETHYLBENZENE	5	U
C245	STYRENE	5	U
C250	TOTAL XYLEMES	5	U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4000

SPECTRIX DC #

SAMPLE NUMBER

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE

QC REPORT NO. : ----

ANALYST: CKT

DATAFILE:

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURI
		54	5
		1	1
		4	4
	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT.	0	0
		STD.	STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

RIC
12/19/86 17:28:00
SAMPLE: TRIP BLANK
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

S S 1 DATA 4U12057V10 #1
CALI: 4U12057V10 #3

SCANS 25 TO 600

SPECTRIX DC # -----14

100.0
RIC
INST ID: 4000

118

TS1

175

SU1

232

TS2

386

SU2

459

TS3

482

SU3

555

58560.

100
5:00

200
10:00

300
15:00

400
20:00

500
25:00

600
30:00 SCAN TIME

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

AMPLE ID. TRIP BLANK

FILENAME 4U12057V10 INST 4000

CLIENT KEYSTONE ANALYST CKT

DATE INJECTED 12/19/86 17:28:00 VERIFIED BY *Not*

CALIB STD DATE 12-14-86 CORR. FACTOR 1.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
176 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000 18271
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000 83931
482 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000 78575
232 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	91 %	1.318 24670
459 CS05	D8-TOLUENE	**SU2**	98	99 %	0.952 73790
555 CS10	F-BROMOFLUOROBENZENE	**SU3**	95	94 %	1.151 50308
118 C030	METHYLENE CHLORIDE *		84	49 UG/L	0.670 32058

* BLANK CORRECTED ON DATA SHEET

CHECKLIST

SECTION III - Blanks

OADS

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram with integrator print out

IADS

004157

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: LLW BLANK

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO : 86-----
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *YPT*

CASE NO. : -----
QC REPORT NO. : -----
CONTRACT NO. : -----
DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW
DATE EXTRACTED/PREPARED: 12/18/86
DATE ANALYZED: 12/18/86

DATAFILE: 4B1218V01
DILUTION FACTOR: 1.00

CAS #	COMPOUND	UG/L	OO
C010	CHLOROMETHANE	10 U	5
C015	BROMOMETHANE	10 U	1
C020	VINYL CHLORIDE	10 U	
C025	CHLOROETHANE	10 U	
C030	METHYLENE CHLORIDE	5	
C035	ACETONE	10 U	
C040	CARBON DISULFIDE	5 U	
C045	1, 1-DICHLOROETHENE	5 U	
C050	1, 1-DICHLOROETHANE	5 U	
C055	TRANS-1, 2-DICHLOROETHENE	5 U	
C060	CHLOROFORM	5 U	
C065	1, 2-DICHLOROETHANE	5 U	
C110	2-BUTANONE	10 U	
C115	1, 1, 1-TRICHLOROETHANE	5 U	
C120	CARBON TETRACHLORIDE	5 U	
C125	VINYL ACETATE	10 U	
C130	BROMODICHLOROMETHANE	5 U	
C140	1, 2-DICHLOROPROPANE	5 U	
C145	TRANS-1, 3-DICHLOROPROPENE	5 U	
C150	TRICHLOROETHENE	5 U	
C155	DIBROMOCHLOROMETHANE	5 U	
C160	1, 1, 2-TRICHLOROETHANE	5 U	
C165	BENZENE	5 U	
C170	CIS-1, 3-DICHLOROPROPENE	5 U	
C175	2-CHLOROETHYL VINYL ETHER	10 U	
C180	BROMOFORM	5 U	
C210	2-HEXANONE	10 U	
C205	4-METHYL-2-PENTANONE	10 U	
C220	TETRACHLOROETHENE	5 U	
C225	1, 1, 2, 2-TETRACHLOROETHANE	5 U	
C230	TOLUENE	5 U	
C235	CHLOROBENZENE	5 U	
C240	ETHYL BENZENE	5 U	
C245	STYRENE	5 U	
C250	TOTAL XYLEMES	5 U	

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: LLW BLK

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 8612057
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/12/86..

SEMICOLATIVES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/18/86..
 DATE ANALYZED: 01/16/87

DATAFILE: 4B12057C01

DILUTION FACTOR: 1:0....

CAS #	COMPOUND	UG/L	004159
C315	PHENOL	10 U	
C325	BIS(2-CHLOROETHYL)ETHER	10 U	
C330	2-CHLOROPHENOL	10 U	
C335	1, 3-DICHLOROBENZENE	10 U	
C340	1, 4-DICHLOROBENZENE	10 U	
C345	BENZYL ALCOHOL	10 U	
C350	1, 2-DICHLOROBENZENE	10 U	
C355	2-METHYLPHENOL	10 U	
C360	BIS(2-CHLOROISOPROPYL)ETHER	10 U	
C365	4-METHYLPHENOL	10 U	
C370	N-NITROSODIPropylAMINE	10 U	
C375	HEXAChLORoETHANE	10 U	
C410	NITROBENZENE	10 U	
C415	ISOPHORONE	10 U	
C420	2-NITROPHENOL	10 U	
C425	2, 4-DIMETHYLPHENOL	10 U	
C430	BENZOIC ACID	50 U	
C435	BIS(2-CHLOROETHOXY)METHANE	10 U	
C440	2, 4-DICHLOROPHENOL	10 U	
C445	1, 2, 4-TRICHLOROBENZENE	10 U	
C450	NAPHTHALENE	10 U	
C455	4-CHLORoANILINE	10 U	
C460	HEXAChLOROBUTADIENE	10 U	
C465	P-CHLORo-M-CRESOL	10 U	
C470	2-METHYLNAPHTHALENE	10 U	
C510	HEXAChLOROCYCLOPENTADIENE	10 U	
C515	2, 4, 6-TRICHLOROPHENOL	10 U	
C520	2, 4, 5-TRICHLOROPHENOL	50 U	
C525	2-CHLORoNAPHTHALENE	10 U	
C530	2-NITROANILINE	50 U	
C535	DIMETHYL PHTHALATE	10 U	
C540	ACENAPHTHYLENE	10 U	
C545	3-NITROANILINE	50 U	
C550	ACENAPHTHENE	10 U	
C555	2, 4-DINITROPHENOL	50 U	
C560	4-NITROPHENOL	50 U	
C565	DIBENZOFURAN	10 U	
C570	2, 4-DINITROTOLUENE	10 U	
C575	2, 6-DINITROTOLUENE	10 U	

SPECTRIX DC # ---- 8

SAMPLE NUMBER: LLW BLK

EMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4812057C01

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	10 U
C585	4-CHLOROPHENYL PHENYL ETHER	10 U
C590	FLUORENE	10 U
C595	4-NITROANILINE	50 U
C610	4, 6-DINITRO-2-METHYLPHENOL	50 U
C615	N-NITROSODIPHENYLAMINE	10 U
C625	4-BROMOPHENYL PHENYL ETHER	10 U
C630	HEXACHLOROBENZENE	10 U
C635	PENTACHLOROPHENOL	50 U
C640	PHENANTHRENE	10 U
C645	ANTHRACENE	10 U
C650	DI-N-BUTYL PHTHALATE	10 U
C655	FLUORANTHENE	10 U
C715	PYRENE	10 U
C720	BUTYL BENZYL PHTHALATE	10 U
C725	3, 3'-DICHLOROBENZIDINE	20 U
C730	BENZO(A)ANTHRACENE	10 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	10 U
C740	CHRYSENE	10 U
C760	DI-N-OCTYL PHTHALATE	10 U
C765	BENZO(B)FLUORANTHENE	10 U
C770	BENZO(K)FLUORANTHENE	10 U
C775	BENZO(A)PYRENE	10 U
C780	INDENO(1, 2, 3-CD)PYRENE	10 U
C785	DIBENZO(A, H)ANTHRACENE	10 U
C790	BENZO(GHI)PERYLENE	10 U

004160

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-057-01A

DC No: KEYSTONE
SMO No: L.BLANK LLW

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/18/86
Date Analyzed: 1/6/87
Dilution Factor: 1
Percent Moisture(decaneted):

GPC Cleanup Yes No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCOLOR-1016	0.50 U
11104-28-2	AROCOLOR-1221	0.50 U
11141-16-5	AROCOLOR-1232	0.50 U
53469-21-9	AROCOLOR-1242	0.50 U
12672-29-6	AROCOLOR-1248	0.50 U
11097-69-1	AROCOLOR-1254	1.0 U
11096-82-5	AROCOLOR-1260	1.0 U

Vs = 500 = Volume of water extracted (ml)

Vt = 5000 = Volume of total extract (ul)

Vi = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 N/A.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: LLW BLANK

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: NOT

DATAFILE: 481218V01

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
4				
0				
1			UC/	
4				
0	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.			
0				
0				

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4930

SPECTRIX DC # --- 8

SAMPLE: LLW BLK

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO. : ---

DC REPORT NO. : ---

ANALYST: QU

DATAFILE: 4B12057C01

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
10-12-3	2-HEXANONE, 5-METHYL-	288	927	M 4G/L
	UNKNOWN--DOESN'T MATCH ANY LIBRARY SPECTRA	1864		412 J
	UNKNOWN--DOESN'T MATCH ANY LIBRARY SPECTRA	1913		021 J 23 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

RIC
12/18/86 9:45:00
SAMPLE: LLW BLANK
COND.: ---
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4B1218U01 #1
CALI: 4B1218U01 #3

SCANS 25 TO 600

SPECTRIX DC # -----14

100.0
RIC
INST ID: 4000

178

SU1

233

122

201

339

365

416

386

TS1

TS2

TS3

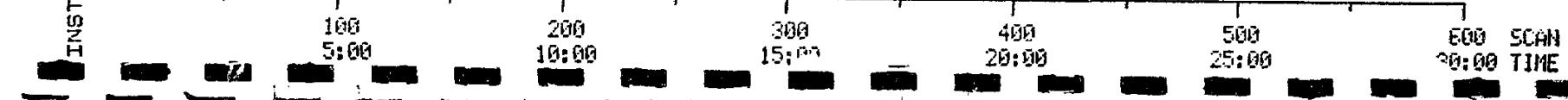
459

482

SU2

556

59200



SPECTRIX DC # -----14

VOA QUANTITATION REPORT

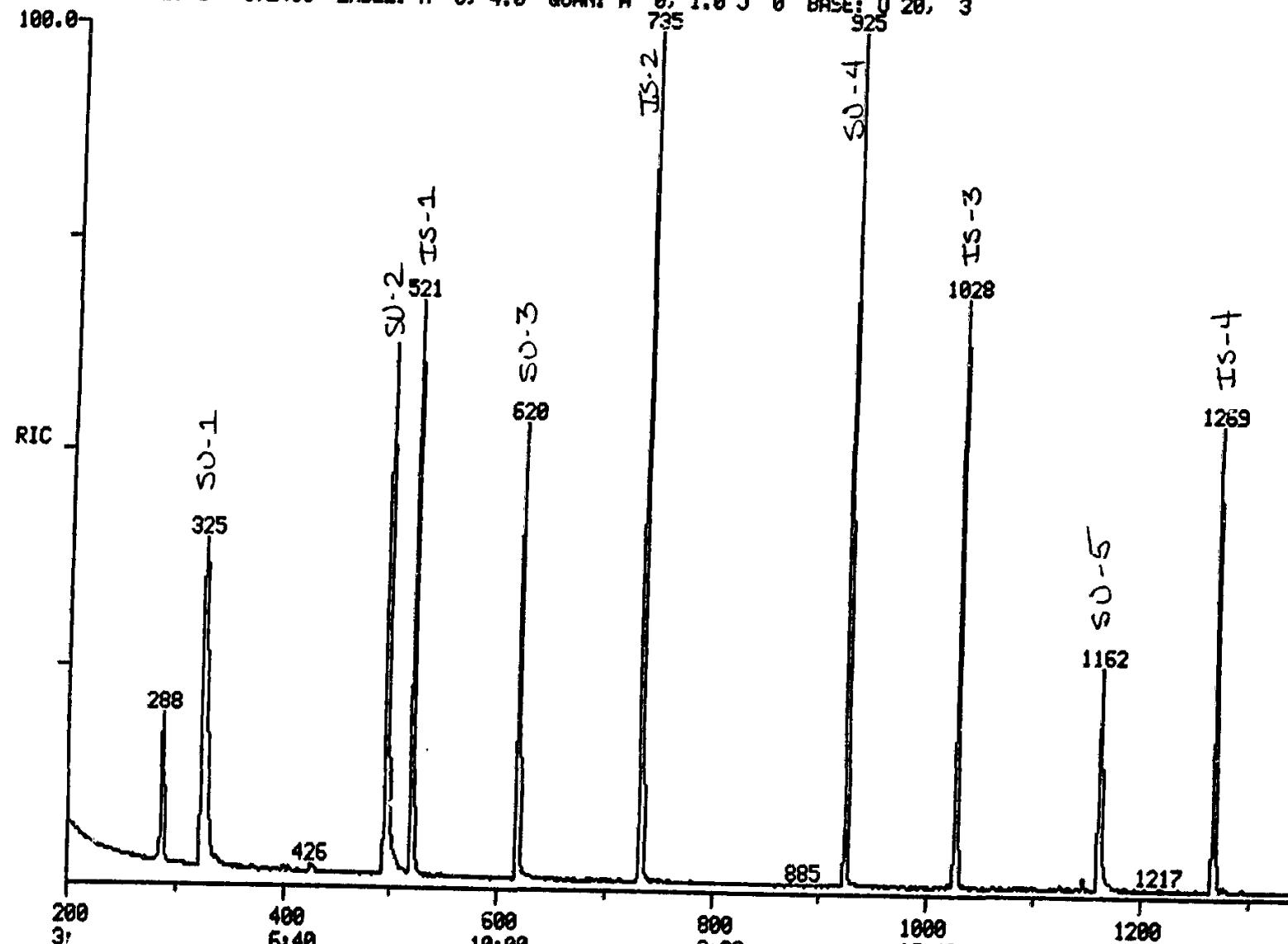
AMPLE ID.	LLW BLANK		
FILENAME	4B121BV01	INST	4000
CLIENT	Keystone	ANALYST	NOT
DATE INJECTED	12/18/86 9:45:00	VERIFIED BY	NOT
CALIB STD DATE	12-14-86	CORR. FACTOR	1.00

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
178 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000 16093
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000 70854
482 CI20	2, 5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000 59304
233 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	93 %	1.309 21452
459 CS05	D8-TOLUENE	**SU2**	98	99 %	0.952 61874
555 CS10	F-BROMOFLUOROBENZENE	**SU3**	95	95 %	1.151 4I461
122 C030	METHYLENE CHLORIDE		84	5 UG/L	0.685 3108
				0	0

INST ID: 4530

SPECTRIX DC # ----14

RIC
01/16/87 14:49:00
SAMPLE: LLW BLK
COND.: ---
RANGE: G 1,2400 LABEL: N 0. 4.0 QUAN: A 0. 1.0 J 0 BASE: U 20, 3
9914166
DATA: 4B12057C01 #1
CALI: 4B12057C01 #3
SCANS 200 TO 1350

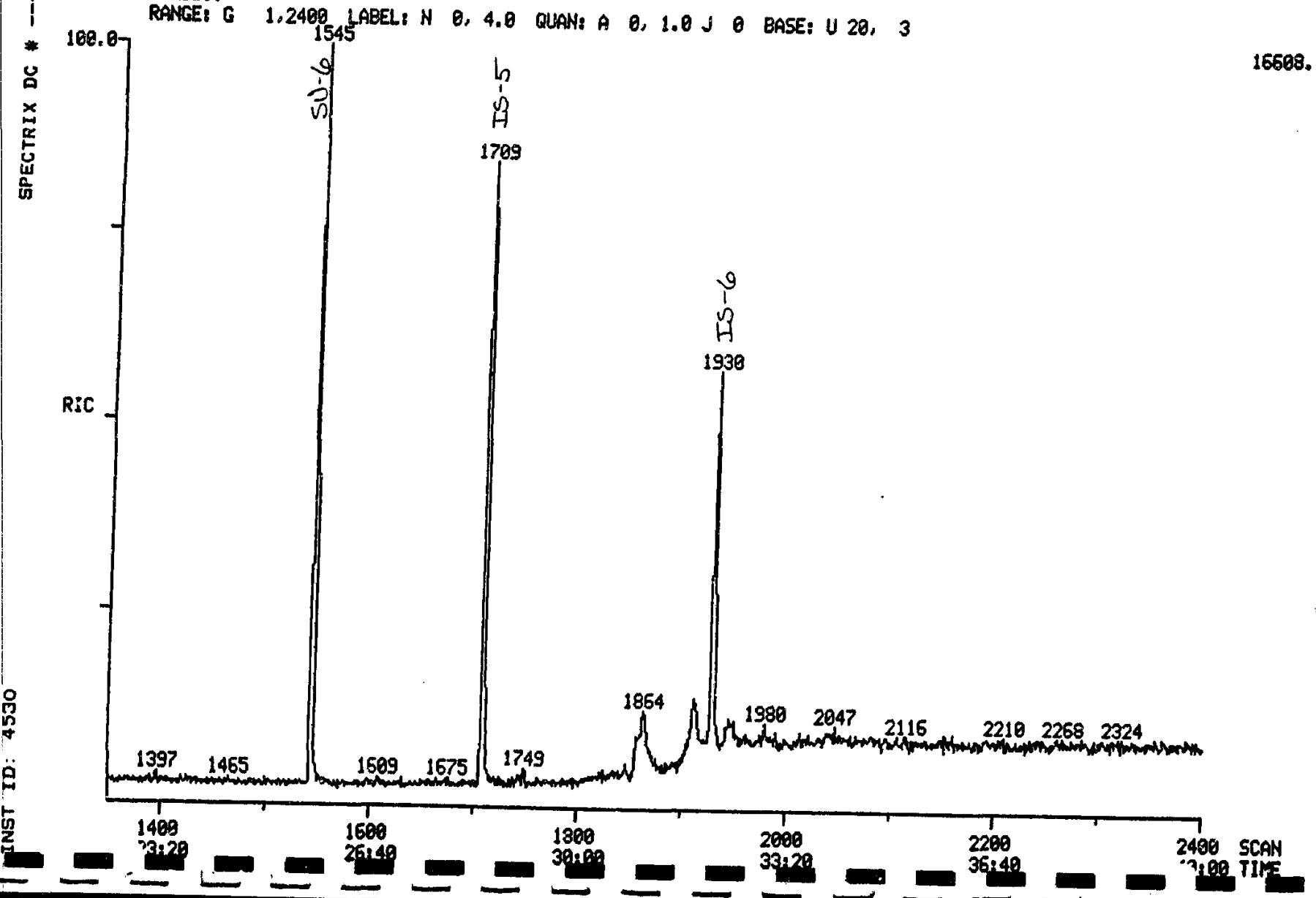


SCAN
TIME

RIC
01/16/87 14:49:00
SAMPLE: LLW BLK
COND.: ---
RANGE: G 1.2400

L 917.00
DATA: 4B12057C01 #1
CALI: 4B12057C01 #3
SCANS 1350 TO 2400

SPECTRIX DC # ----14



SPECTRIX DC # -----14

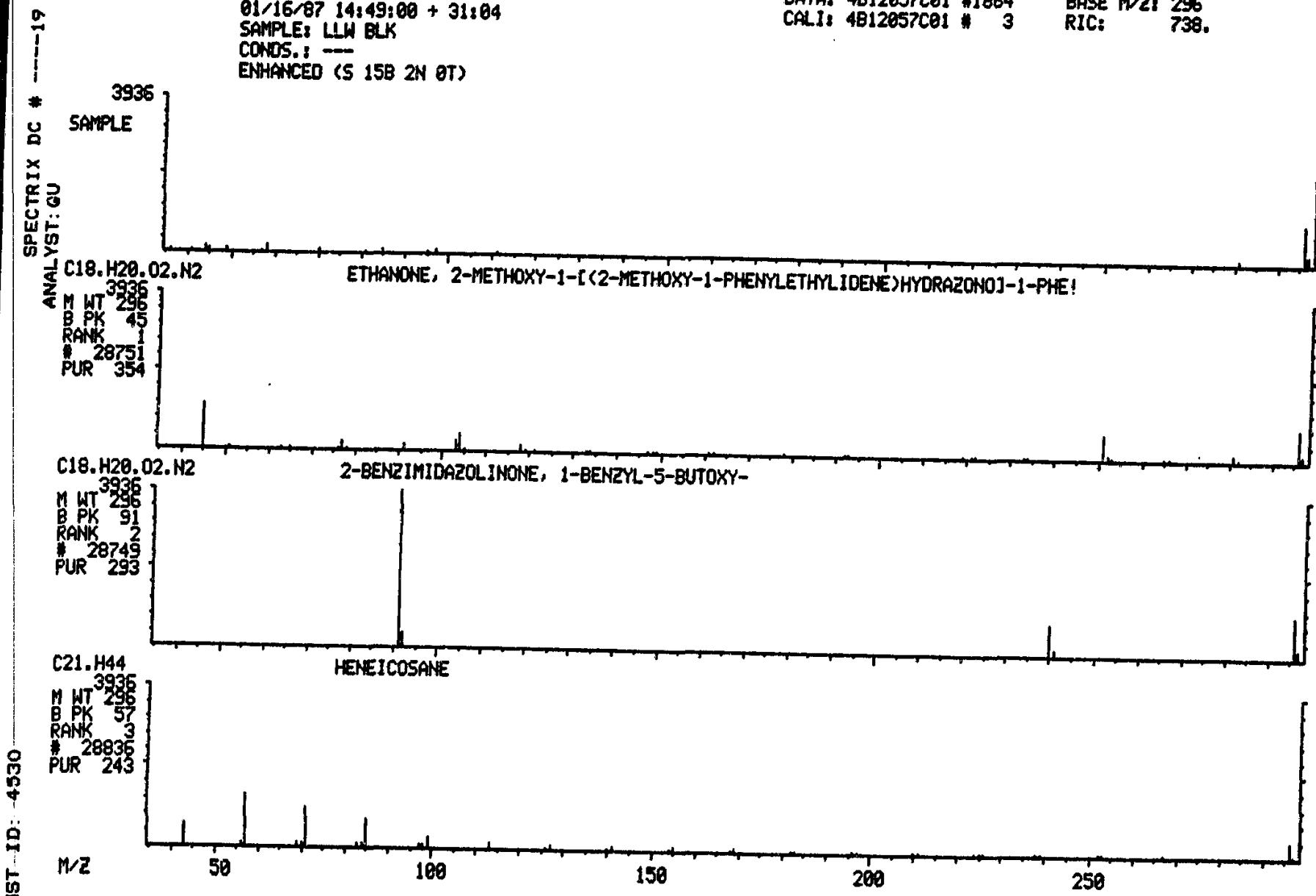
SEMIVOLATILE WATER REPORT

SAMPLE ID.	LLW BLK		
ILENAME	4B12057C01	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/16/87 14:49:00	VERIFIED BY	<i>JHH</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	1.00

SCAN#		SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
521	CI30	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	40 UG/L	1.000	1435:
735	CI40	NAPHTHALENE-D8 ** IS2 **	136	40 UG/L	1.000	5598:
1028	CI50	ACENAPHTHENE-D10 ** IS3 **	164	40 UG/L	1.000	2756:
1269	CI60	PHENANTHRENE-D10 ** IS4 **	188	40 UG/L	1.000	03430:
1709	CI70	CHRYSENE-D12 ** IS5 **	240	40 UG/L	1.000	1854:
1930	CI75	PERYLENE-D12 ** IS6 **	264	40 UG/L	1.000	1321:
325	CS50	2-FLUOROPHENOL ** SU1 **	112	90 %	0.624	3104:
497	CS45	PHENOL-D5 ** SU2 **	99	67 %	0.954	4777:
1162	CS55	2,4,6-TRIBROMOPHENOL ** SU5 **	330	65 %	1.130	0 491:
620	CS20	NITROBENZENE-D5 ** SU3 **	82	72 %	0.844	C2440:
925	CS25	2-FLUOROBIPHENYL ** SU4 **	172	68 %	0.900	4915:
1545	CS30	TERPHENYL-D14 ** SU6 **	244	79 %	0.904	1856:

LIBRARY SEARCH 691700
01/16/87 14:49:00 + 31:04
SAMPLE: LLW BLK
CONDNS.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4B12057C01 #1864
CALI: 4B12057C01 # 3
BASE M/Z: 296
RIC: 738.



INST ID: 4530

004169

SPECTRUM DC # ---19

LIBRARY SEARCH 004170
01/16/87 14:49:00 + 31:53
SAMPLE: LLW BLK
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4B12057C01 #1913 BASE M/Z: 85
CALI: 4B12057C01 # 3 RIC: 940.



M WT 1523
B PK 312
RANK 55
30354
PUR 422

M WT 1523
B PK 289
RANK 67
271162
PUR 4147

M WT 1523
B PK 248
RANK 57
224453
PUR 405

INST ID: 4530

M/Z 50 100 150 200 250 300

INST ID 4000

SPECTRIX DC # ---- 8

SAMPLE NUMBER: LLW BLANK

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX

CASE NO. : ----

LAB SAMPLE ID NO. : 86----

QC REPORT NO. : ----

SAMPLE MATRIX: WATER

CONTRACT NO. : ----

DATA RELEASE AUTHORIZED BY: *ypm*

DATE SAMPLE RECEIVED: 12/12/86

VOLATILES

CONCENTRATION: LOW

DATAFILE: 4B1219V01A

DATE EXTRACTED/PREPARED: 12/19/86

DILUTION FACTOR: 1.00

DATE ANALYZED: 12/19/86

CAS #	COMPOUND	UG/L	7
C010	CHLOROMETHANE	10 U	4
C015	BROMOMETHANE	10 U	0
C020	VINYL CHLORIDE	10 U	0
C025	CHLOROETHANE	10 U	0
C030	METHYLENE CHLORIDE	10 U	0
C035	ACETONE	5	0
C040	CARBON DISULFIDE	10 U	0
C045	1, 1-DICHLOROETHENE	5 U	0
C050	1, 1-DICHLOROETHANE	5 U	0
C055	TRANS-1, 2-DICHLOROETHENE	5 U	0
C060	CHLOROFORM	5 U	0
C065	1, 2-DICHLOROETHANE	5 U	0
C110	2-BUTANONE	5 U	0
C115	1, 1, 1-TRICHLOROETHANE	10 U	0
C120	CARBON TETRACHLORIDE	5 U	0
C125	VINYL ACETATE	5 U	0
C130	BROMODICHLOROMETHANE	10 U	0
C140	1, 2-DICHLOROPROPANE	5 U	0
C145	TRANS-1, 3-DICHLOROPROPENE	5 U	0
C150	TRICHLOROETHENE	5 U	0
C155	DIBROMOCHLOROMETHANE	5 U	0
C160	1, 1, 2-TRICHLOROETHANE	5 U	0
C165	BENZENE	5 U	0
C170	CIS-1, 3-DICHLOROPROPENE	5 U	0
C175	2-CHLOROETHYL VINYL ETHER	5 U	0
C180	BROMOFORM	10 U	0
C210	2-HEXANONE	5 U	0
C205	4-METHYL-2-PENTANONE	10 U	0
C220	TETRACHLOROETHENE	10 U	0
C225	1, 1, 2, 2-TETRACHLOROETHANE	5 U	0
C230	TOLUENE	5 U	0
C235	CHLOROBENZENE	5 U	0
C240	ETHYLBENZENE	5 U	0
C245	STYRENE	5 U	0
C250	TOTAL XYLEMES	5 U	0

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4000

SPECTRIX DG # ----- 8

SAMPLE NUMBER: LLW BLANK

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO. : -----

QC REPORT NO. : -----

ANALYST: NOT

DATAFILE: 4B1219V01A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
		4		
		17	UG/	
	NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.	4	0	0

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

RIC
12/19/85 10:00:00
SAMPLE: LLW BLANK
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

C L 1 4B1213U01A #1
CALI: 4B1213U01A #3
SCANS 25 TO 600

54464.

SPECTRIX DC # -----14

100.0

RIC

INST ID 4000

1

1

1

1

TS1

179

SU1

233

123

142

202

TS2

366

SU2

459

TS3

482

SU3

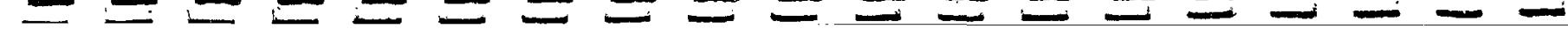
556

600
30:00 SCAN TIME

15:00

10:00

5:00



004173

SPECTRIX DC # -----14

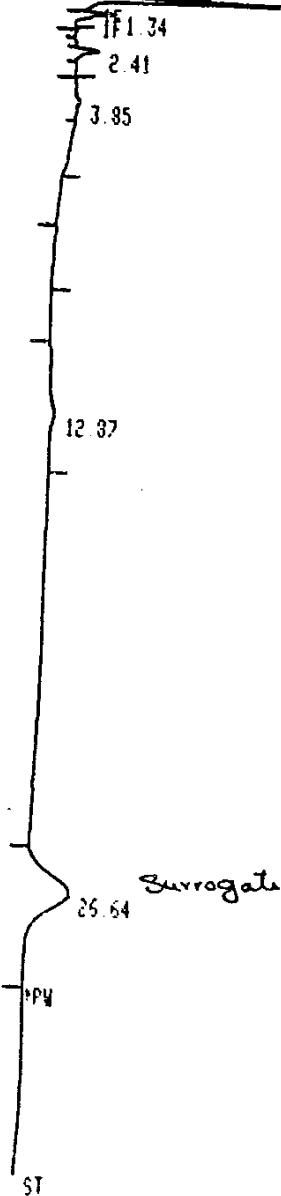
VOA QUANTITATION REPORT

AMPLE ID.	LLW BLANK				
FILENAME	481219V01A	INST	4000		
CLIENT	KEYSTONE	ANALYST	NOT		
DATE INJECTED	12/19/86 10:00:00	VERIFIED BY	NOT		
CALIB STD DATE	12-14-86	CORR. FACTOR	1. 00		

SCAN#	VOA COMPOUNDS	M/E	AMOUNT	RT	AREA
179 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1. 000 14734
386 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1. 000 63765
482 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1. 000 52138
233 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	90 %	1. 302 49726
459 CS05	DB-TOLUENE	**SU2**	98	102 %	0. 952 54633
555 CS10	F-BROMOFLUOROBENZENE	**SU3**	95	95 %	1. 151 36393
123 C030	METHYLENE CHLORIDE		84	5 UG/L	0. 687 2909

STAK-LF 86-12-057-01A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. Lab. Blank (L)
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 μL
INSTRUMENT ID. 6600

RUN # 18
WORKFILE ID: C
WORKFILE NAME:

JAN/06/87 21:27:04

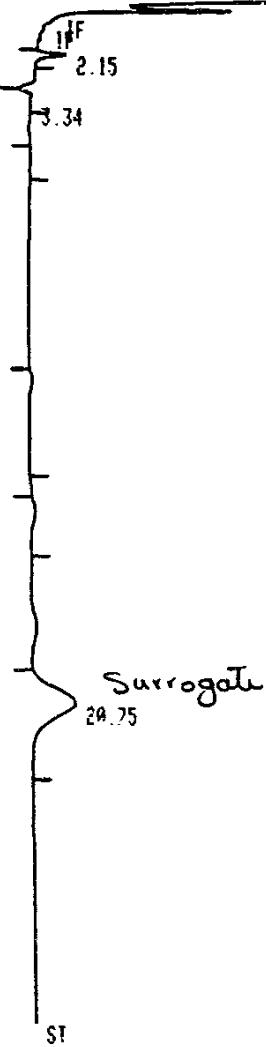
HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.34	7927	D BB	0.089	27.958
2.41	5917	VV	0.140	20.869
3.85	2228	BV	0.664	7.858
12.87	1354	PB	0.997	4.811
26.64	10917	BB	1.165	38.584

TOTAL HGT= 28353
MUL FACTOR= 1.0000E+00

START #: 56-012-032-01A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. KeyStone

SAMPLE NO. L. Blank (u)
PESTICIDE FRACTION
1ST COLUMN QUANT.

1ST COLUMN ID
1.8m x 4mm ID x6mm (c)
1.5% SP 2250
1.95% SP 2401
100/120 SUPELCOPORT

COLUMN TEMP. 203°C
VOLUME INJ. 1µl
INSTRUMENT ID. 5600C

00

RUN #: 181 JAN/12/87 13:39:58
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%				
RT	HEIGHT	TYPE	AR/HT	HEIGHT%
2.15	3714	BB	0.133	35.517
3.34	1279	PB	0.297	12.231
20.75	5464	BB	0.966	52.252

TOTAL HGT= 10457
MUL FACTOR= 1.0000E+00

Page 1

Received: 12/12/86

SPECTRIX

REPORT

WORK ORDER # 00-12-03/

LL14700 Results by Sample

SAMPLE ID LAB BLANK

SAMPLE # 01 FRACTIONS: A

Date & Time Collected 12/12/86

Category

AG	<10	AS	<10	BE	<3	CD	<5	CR	<9	CU	<10
	ug/L										
FE	<30	HG	<0.2	NI	<29	PB	<5	SB	<44	SE	<5
	ug/L										
TL	<10	ZN	<8								
	ug/L		ug/L								

GC/MS Tuning and Mass Calibration

DECAFLUOROTRIPHENYLPHOSPHINE

Case Number:

Inst ID:

4530

Laboratory:

Lab ID:

ES0112C01

Sens Date: 01/12/87

Cali Date:

Contract:

Sens Time:

9:32:00

Analyst: GU

Data release authorized by:

M/E

Ion Abundance Criteria

Spec # 390

51	30 to 60% of mass 198	
68	less than 2% of mass 69	46.34
69	mass 69 relative abundance	0.00 (0.00) 1
70	less than 2% of mass 69	22.42
127	40 TO 60% of mass 198	0.12 (0.56) 1
197	less than 1% of mass 198	55.38
198	base peak, 100% relative abundance	0.00
199	5 TO 9% of mass 198	100.00
275	10 TO 30% of mass 198	6.34
363	greater than 1% of mass 198	21.52
441	less than mass 443	1.29
442	greater than 40% of mass 198	9.09
443	17 TO 23% of mass 442	75.61
		14.47 (19.14) 2

1 - Value in parenthesis is % of mass 69

2 - Value in parenthesis is % of mass 442

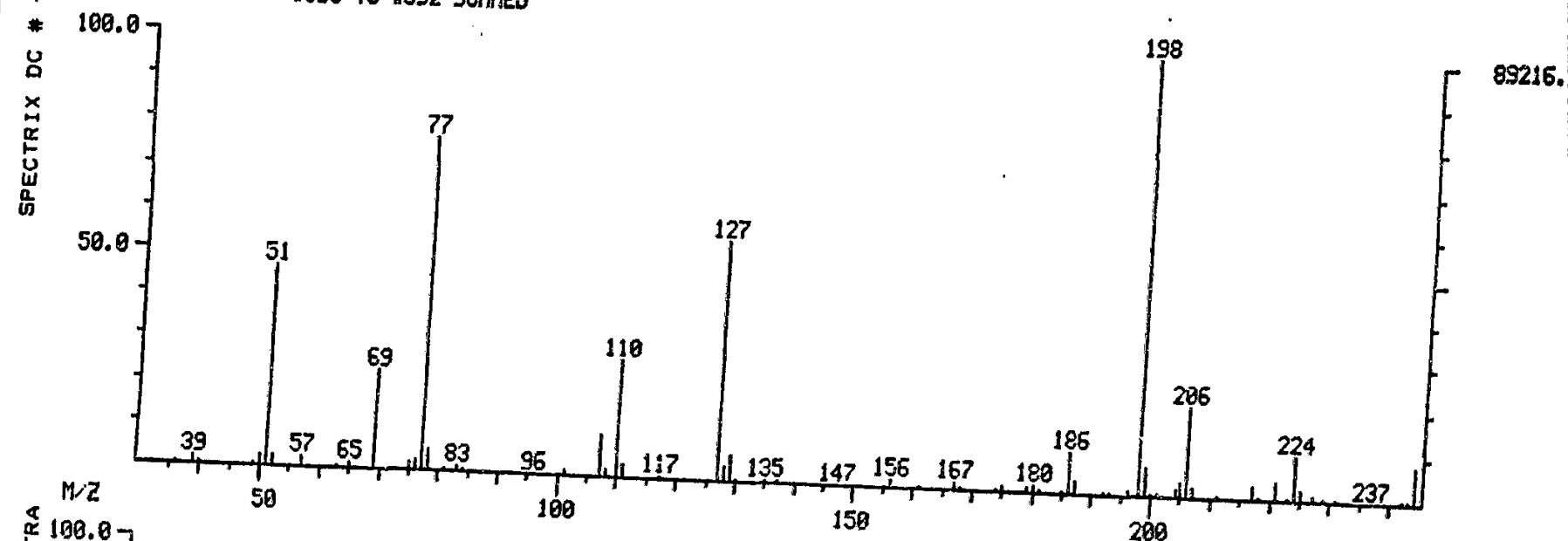
004178

SPECTRUM DC # ---15

MASS SPECTRUM
01/12/87 9:32:00 + 6 L 1 7 0 0
SAMPLE: DFTPP
COND.: ---
#390 TO #392 SUMMED

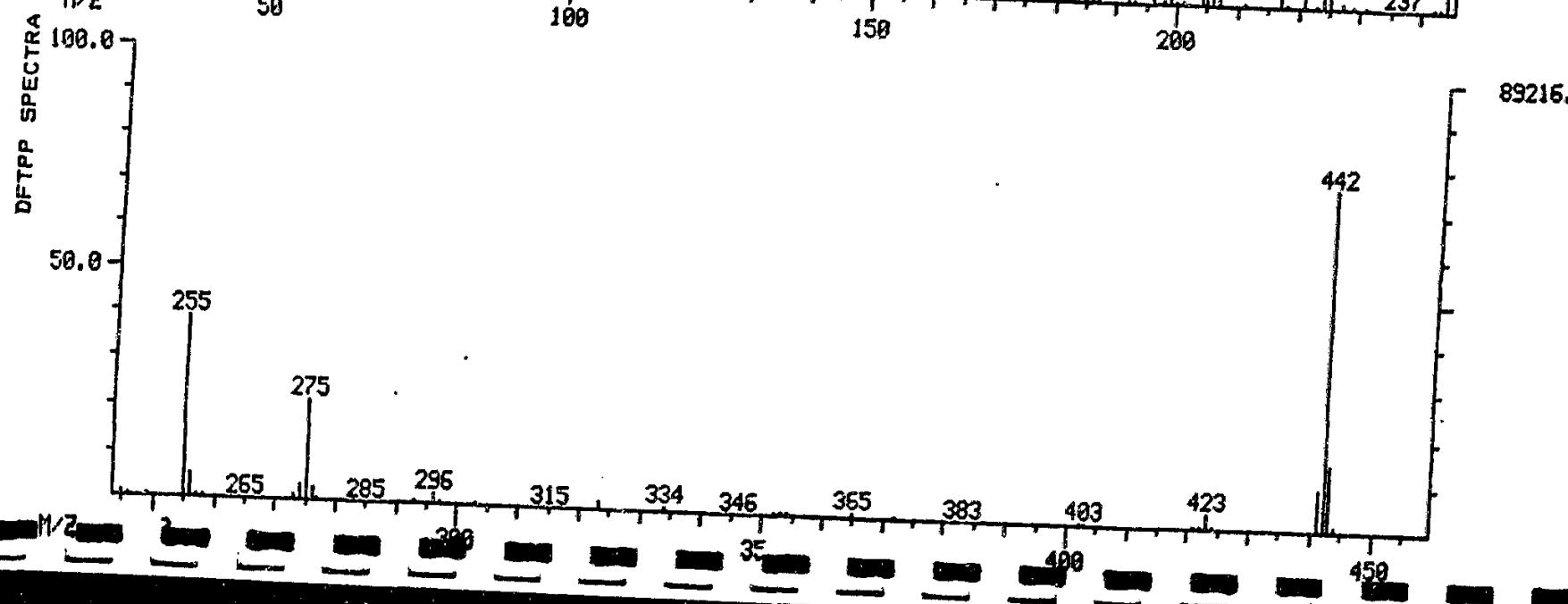
DATA: ES0112C01 #391
CALI: ES0112C01 #3

BASE M/Z: 198
RIC: 598016.



DFTPP SPECTRA

INST ID: 4530



INST ID: 4530

SPECTRIX DC # ----15

Mass List

01/12/87 9:32:00 + 6:31

Data: ES0112C01 # 391

Base m/z: 198

Cali: ES0112C01 # 3

RIC: 598016.

Sample: DFTPP

Tonds.: ---

390 to #392 summed

Mass	% RA	Inten.	Minima	Min	Inten:	250.
			Maxima	#	O	
36	0.00	0.				
444						
36?	0.66	586.	204	1.	59	1422.
39?	1.72	1536.	205	3.	73	3328.
40	0.56	500.	206	20.	09	17920.
49	0.68	611.	207	2.	47	2204.
50	2.37	2116.	211	0.	72	639.
51	46.34	41344.	217	2.	75	2452.
52	2.20	1962.	218	0.	43	382.
57	2.11	1880.	221	4.	28	3820.
63	0.33	298.	223	0.	85	761.
65	1.31	1170.	224	9.	77	8720.
69	22.42	20000.	225	2.	37	2112.
75	1.56	1390.	227	1.	23	1096.
76	2.18	1946.	229	0.	60	539.
77	76.61	68352.	231	0.	39	346.
78	4.77	4256.	242	0.	42	374.
81	0.59	528.	243	0.	45	400.
83	1.10	983.	244	8.	13	7256.
84	0.39	349.	245	0.	99	887.
101	0.98	878.	246	0.	48	427.
107	9.42	8400.	249	0.	29	263.
108	1.97	1762.	255	38.	74	34560.
110	27.01	24096.	256	5.	51	4912.
111	2.63	2344.	257	0.	35	312.
117	0.32	288.	258	0.	73	649.
127	55.38	49408.	273	1.	24	1106.
128	3.66	3264.	274	3.	70	3300.
129	5.94	5296.	275	21.	52	19200.
130	0.42	374.	276	2.	68	2388.
135	0.52	460.	277	0.	84	746.
137	0.32	284.	293	0.	30	268.
155	0.40	355.	296	2.	56	2288.
156	1.50	1338.	297	0.	32	285.
161	0.54	483.	303	0.	40	361.
167	1.84	1638.	323	1.	58	1410.
168	0.62	555.	334	0.	88	787.
174	0.36	324.	352	0.	42	375.
175	1.18	1052.	353	0.	32	283.
179	0.94	843.	354	0.	58	516.
180	1.49	1330.	365	1.	29	1152.
181	0.75	671.	372	0.	84	753.
185	0.52	468.	402	0.	31	279.
186	9.11	8128.	403	0.	49	440.
187	2.71	2416.	421	0.	42	373.
192	0.34	305.	422	0.	39	351.
193	0.35	494.	423	3.	30	2940.
196	1.35	1204.	424	0.	59	522.
198	100.00	89216.	441	9.	09	8112.
199	6.34	5656.	442	75.	61	67436.
200	0.39	346.	443	14.	47	12912.
201	0.29	261.	444	1.	29	1148.

004180

GC/MS Tuning and Mass Calibration

DECAFLUOROTRIPHENYLPHOSPHINE

ase Number: inst ID: 4530 Lab ID: ES0113C01	Laboratory: Sens Date: 01/13/87 Cali Date:	Contract: Sens Time: 8:37:00 Analyst: GU
---	--	--

Data release authorized by:

УРН

Contract:
Sens Time: 8:37:00
Analyst: GU

M/E	Ion Abundance Criteria	Spec # 399
51	30 to 60% of mass 198	48. 42
68	less than 2% of mass 69	0. 00
69	mass 69 relative abundance	23. 20
70	less than 2% of mass 69	0. 16
127	40 TO 60% of mass 198	57. 85
197	less than 1% of mass 198	0. 00
198	base peak, 100% relative abundance	100. 00
199	5 TO 9% of mass 198	6. 34
275	10 TO 30% of mass 198	20. 35
365	greater than 1% of mass 198	1. 34
441	less than mass 443	9. 38
442	greater than 40% of mass 198	78. 61
443	17 TO 23% of mass 442	14. 75

1 - value in parenthesis is % of mass 69

2 - value in parenthesis is % of mass 442

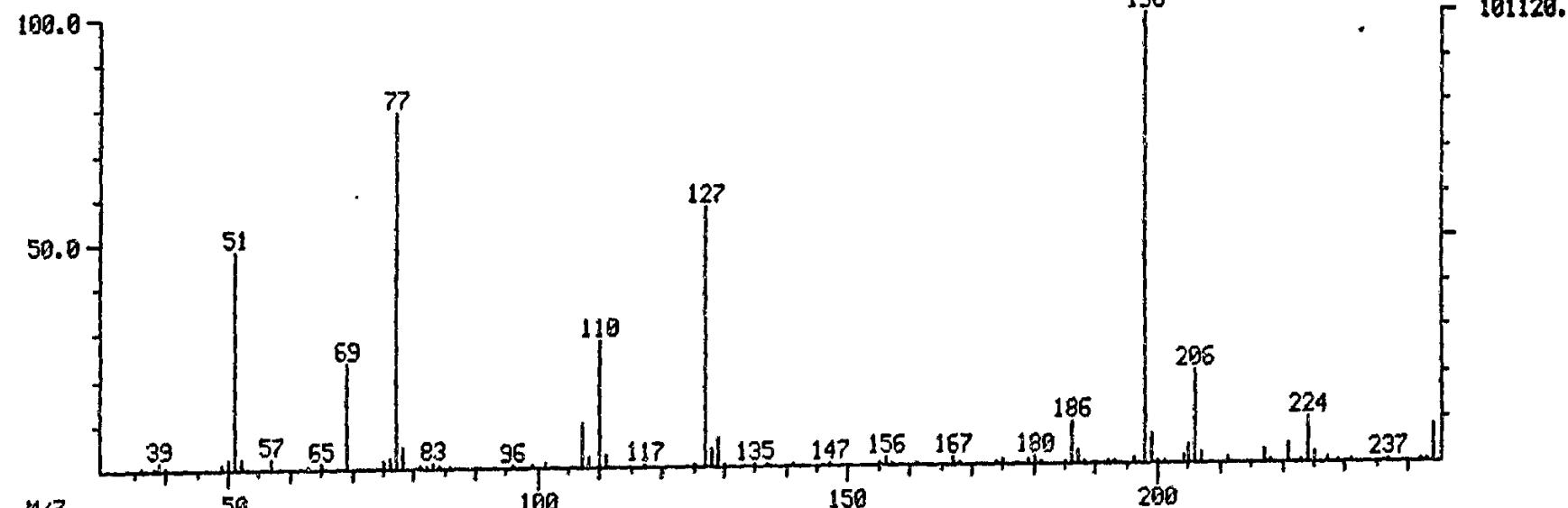
004181

SPECTRIX DC # -----15

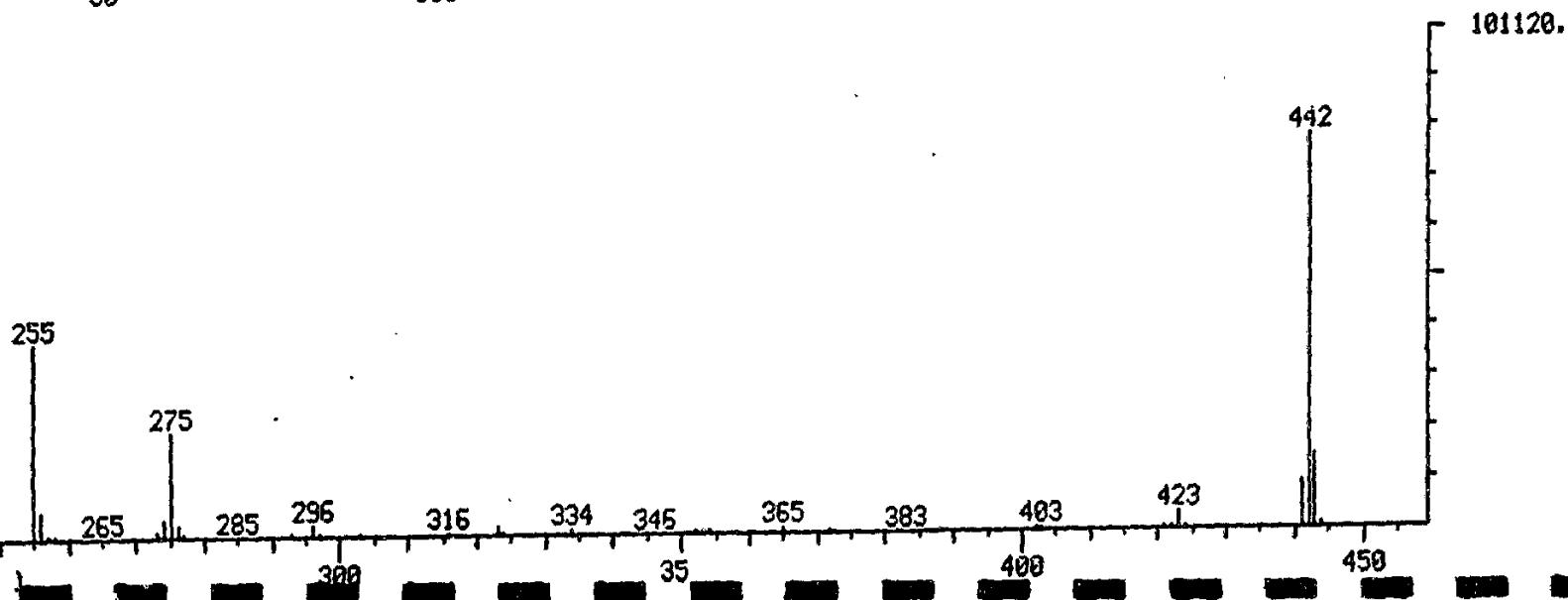
MASS SPECTRUM
01/13/87 8:37:00 + 281 0040
SAMPLE: DFTPP
COND.: ---
#399 TO #402 SUMMED

DATA: E50113C01 #400
CALI: E50113C01 #3

BASE M/Z: 198
RIC: 695296.



DFTPP SPECTRA
INST ID: 4530



INST ID: 4530

DFTPP MASS LISTING

SPECTRIX DC # ----15

Mass List

01/13/87 8:37:00 + 6:40

Data: ES0113C01 # 400

Base m/z: 198

Sample: DFTPP

Calib: ES0113C01 # 3

RIC: 695296.

Bonds: ---

399 to #402 summed

Mass	0.00	0.	Minima Maxima	Min Inten: # 0	340.
	% RA	Inten.	Mass	% RA	Inten.
36?	0.81	821.	200	0.44	440.
39?	1.82	1842.	201	0.38	387.
40	0.70	704.	204	1.68	1702.
49	1.16	1172.	205	3.78	3824.
50	2.56	2584.	206	20.22	20448.
51	48.42	48960.	207	2.53	2556.
52	2.31	2340.	211	0.90	912.
57	2.25	2272.	217	2.85	2884.
63	0.42	424.	218	0.49	491.
65	1.43	1448.	221	4.29	4336.
69	23.20	23455.	223	0.86	667.
75	1.65	1666.	224	9.70	9808.
76	2.31	2332.	225	2.43	2456.
77	79.11	80000.	227	1.26	1278.
78	4.93	4984.	229	0.67	674.
81	0.68	688.	231	0.46	470.
83	1.23	1240.	242	0.47	480.
84	0.69	701.	243	0.49	500.
86	0.47	474.	244	7.94	8032.
101	1.12	1134.	245	1.01	1026.
107	9.86	9968.	246	0.52	526.
108	2.14	2160.	249	0.37	374.
110	27.94	28256.	255	37.59	38016.
111	2.88	2912.	256	5.35	5408.
117	0.41	412.	257	0.39	390.
127	57.85	58496.	258	0.76	767.
128	3.79	3832.	273	1.21	1220.
129	6.16	6232.	274	3.56	3604.
130	0.48	489.	275	20.35	20576.
135	0.60	606.	276	2.57	2600.
137	0.38	386.	277	0.84	854.
141	0.35	358.	293	0.34	340.
155	0.46	467.	296	2.51	2540.
156	1.61	1630.	303	0.47	475.
161	0.62	625.	323	1.52	1536.
167	1.90	1924.	334	0.89	905.
168	0.69	693.	352	0.45	456.
174	0.43	439.	354	0.60	603.
175	1.24	1250.	365	1.34	1350.
179	1.09	1104.	372	0.85	856.
180	1.58	1600.	402	0.36	364.
181	0.80	810.	403	0.50	503.
185	0.60	606.	421	0.52	522.
186	9.30	9408.	422	0.43	433.
187	2.76	2792.	423	3.39	3424.
192	0.41	414.	424	0.65	661.
193	0.62	623.	441	9.38	9488.
196	1.44	1458.	442	78.61	79488.
198	100.00	101120.	443	14.75	14912.
199	6.34	6416.	444	1.36	1372.

004183

GC/MS Tuning and Mass Calibration

DECAFLUOROTRIPHENYLPHOSPHINE

Case Number:

Inst ID: 4530
Lab ID: ES0116C01

Laboratory:

Sens Date: 01/16/87
Cali Date:

Contract:

Sens Time: 10:12:00
Analyst: GU

Data release authorized by:

JPM

M/E	Ion Abundance Criteria	Spec # 411
51	30 to 60% of mass 198	46.44
68	less than 2% of mass 69	0.05 (0.23) 1
69	mass 69 relative abundance	22.60
70	less than 2% of mass 69	0.16 (0.70) 1
127	40 TO 60% of mass 198	59.73
197	less than 1% of mass 198	0.00
198	base peak, 100% relative abundance	100.00
199	5 TO 9% of mass 198	6.19
275	10 TO 30% of mass 198	19.76
365	greater than 1% of mass 198	1.29
441	less than mass 443	9.88
442	greater than 40% of mass 198	80.82
443	17 TO 23% of mass 442	15.50 (19.17) 2

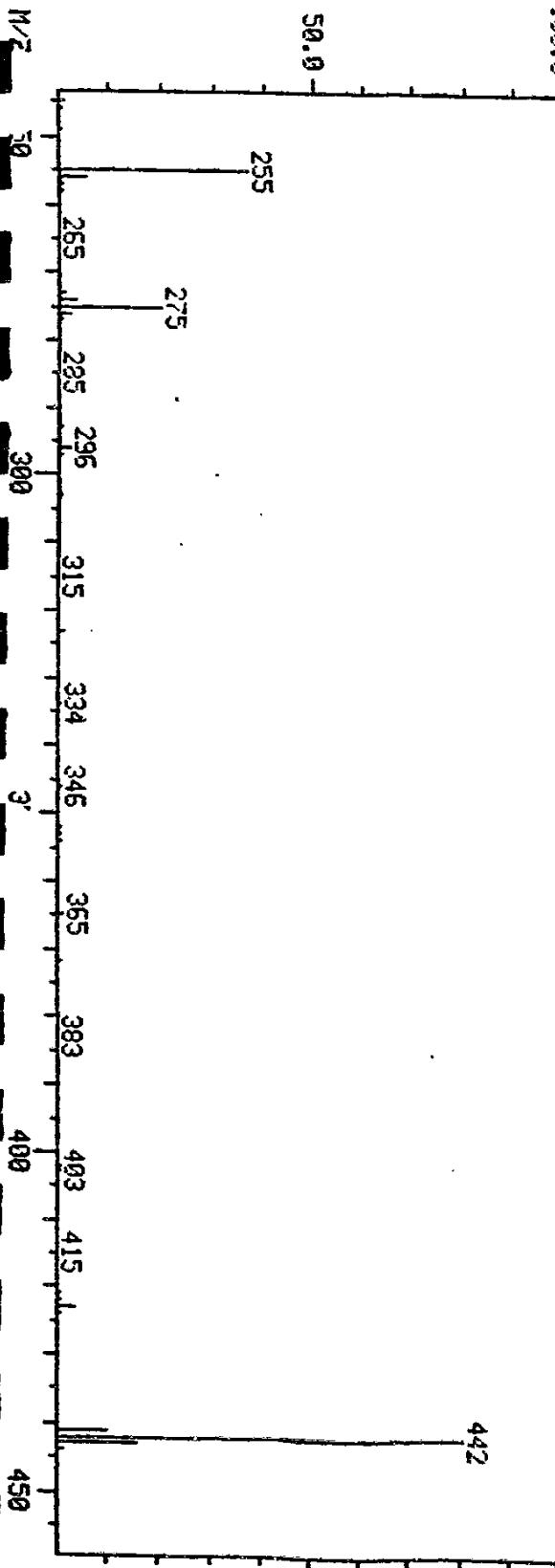
1 - value in parenthesis is % of mass 69

2 - value in parenthesis is % of mass 442

004104

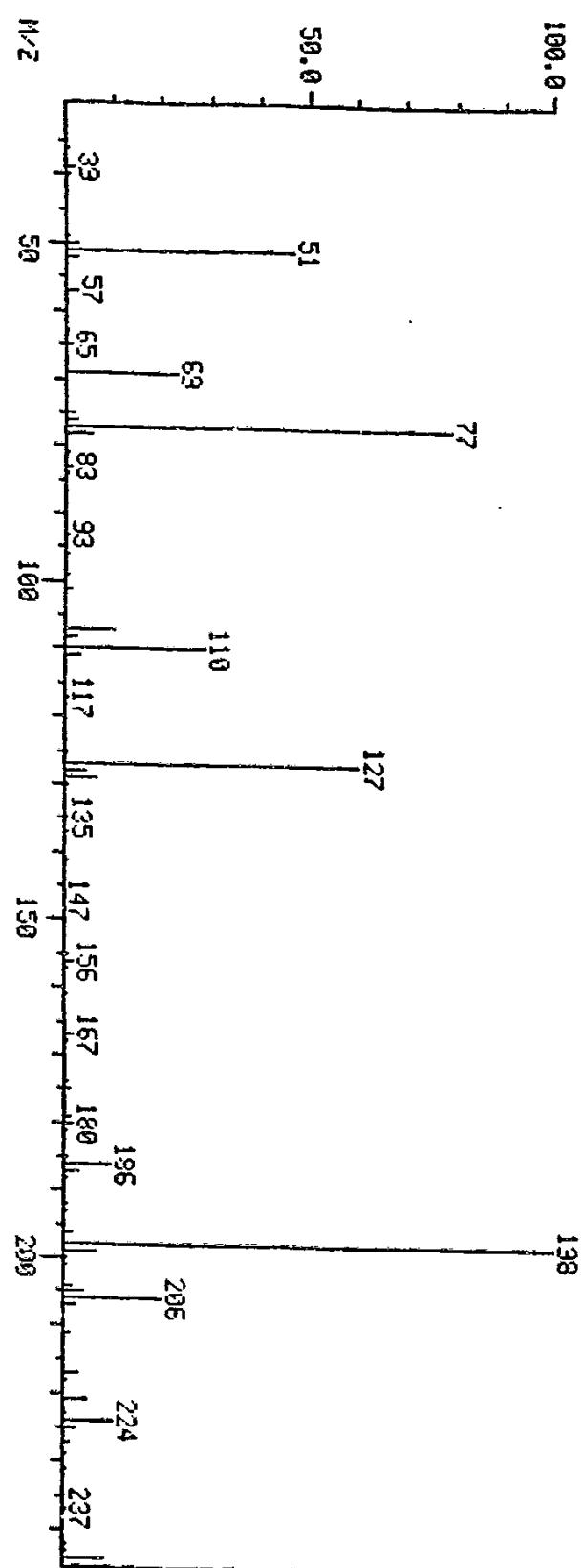
INST ID 4530

DFTPP SPECTRA



SPECTRIX DC # ----15

MASS SPECTRUM
01/16/87 10:12:00 + 5:32
SAMPLE: DFTPP
COND.: ---
#411 TO #413 SUMMED



DATA: ES0116C01 #412
CALL: ES0116C01 #3

BASE M/Z: 198
RIC: 636928.

Mass List

01/16/87 10:12:00 + 6:52
 Sample: DFTPP
 Conds.: ---
 #411 to #413 summed

Data: ES0116C01 # 412
 Cali: ES0116C01 # 3

Base m/z: 196
 RIC: 636928.

Mass	36	0.00	0.	Minima	Min Inten:	
	445	% RA	Inten.	Maxima	#	321.
	Mass			Mass	0	
	36?	0.63	590.	201	0.38	357.
	39?	1.75	1636.	204	1.62	1512.
	40	0.62	582.	205	3.94	3684.
	49	0.54	505.	206	20.03	18720.
	50	2.29	2140.	207	2.45	2292.
	51	46.44	43392.	211	0.88	821.
	52	2.20	2060.	217	2.92	2732.
	57	2.16	2014.	218	0.45	424.
	63	0.38	357.	221	4.50	4208.
	65	1.42	1324.	223	0.86	807.
	69	22.60	21120.	224	9.64	9008.
	75	1.65	1542.	225	2.41	2248.
	76	2.18	2034.	227	1.26	1174.
	77	78.63	73472.	229	0.67	625.
	78	5.06	4728.	231	0.45	421.
	81	0.64	599.	237	0.36	338.
	83	1.21	1134.	242	0.44	412.
	84	0.38	351.	243	0.51	481.
	101	1.08	1006.	244	7.85	7336.
	107	9.95	9296.	245	0.98	914.
	108	2.17	2032.	246	0.50	466.
	110	28.36	26496.	249	0.38	353.
	111	2.87	2680.	255	36.71	34304.
	117	0.43	404.	256	5.04	4712.
	127	59.73	55808.	257	0.36	340.
	128	3.78	3536.	258	0.73	679.
	129	6.40	5984.	273	1.16	1080.
	130	0.49	459.	274	3.33	3108.
	135	0.57	533.	275	19.76	18464.
	137	0.36	336.	276	2.43	2272.
	141	0.41	380.	277	0.80	751.
	155	0.49	461.	296	2.43	2272.
	156	1.61	1500.	297	0.36	339.
	161	0.62	576.	303	0.45	424.
	167	1.94	1814.	323	1.42	1328.
	168	0.66	619.	334	0.82	767.
	174	0.42	389.	352	0.43	398.
	175	1.27	1190.	354	0.52	488.
	179	1.04	971.	365	1.29	1202.
	180	1.63	1526.	372	0.84	781.
	181	0.81	757.	402	0.34	321.
	185	0.35	518.	403	0.51	475.
	186	9.16	8560.	421	0.48	448.
	187	2.71	2536.	422	0.38	354.
	192	0.42	396.	423	3.42	3200.
	193	0.63	588.	424	0.64	598.
	196	1.52	1418.	441	9.88	9232.
	198	100.00	93440.	442	80.82	75520.
	199	6.19	5784.	443	15.50	14480.
	200	0.43	401.	444	1.44	1350.

GC/MS Tuning and Mass Calibration

BROMOFLUOROBENZENE

File Number:

1st ID: 4000
Lab ID: ES1218V01

Laboratory:

Sens Date: 12/18/86
Cali Date:

Contract:
Sens Time: 8:01:00
Analyst: NOT

Data release authorized by:

[Signature]

m/z

Ion Abundance Criteria

Spec # 207

50	15 to 40% of mass 95	32.75
75	30 to 60% of mass 95	43.08
95	base peak, 100% relative abundance	100.00
96	5 to 9% of mass 95	7.25
173	less than 1% of mass 95	0.00
174	greater than 50% of mass 95	84.40
175	5 to 9% of mass 174	6.91 (8.19) 1
176	between 95% and 101% of mass 174	80.99 (95.96) 1
177	5 to 9% of mass 176	5.59 (6.90) 2

1 - value in parenthesis is % of mass 174

2 - value in parenthesis is % of mass 176

7
8
1
4
0

INST ID: 4000

Mass List

12/18/86 8:01:00 + 10:21

Sample: BFB

Ends: ----

BFB MASS LISTING

SPECTRIX DC # -----15

Data: ES1218V01 # 207

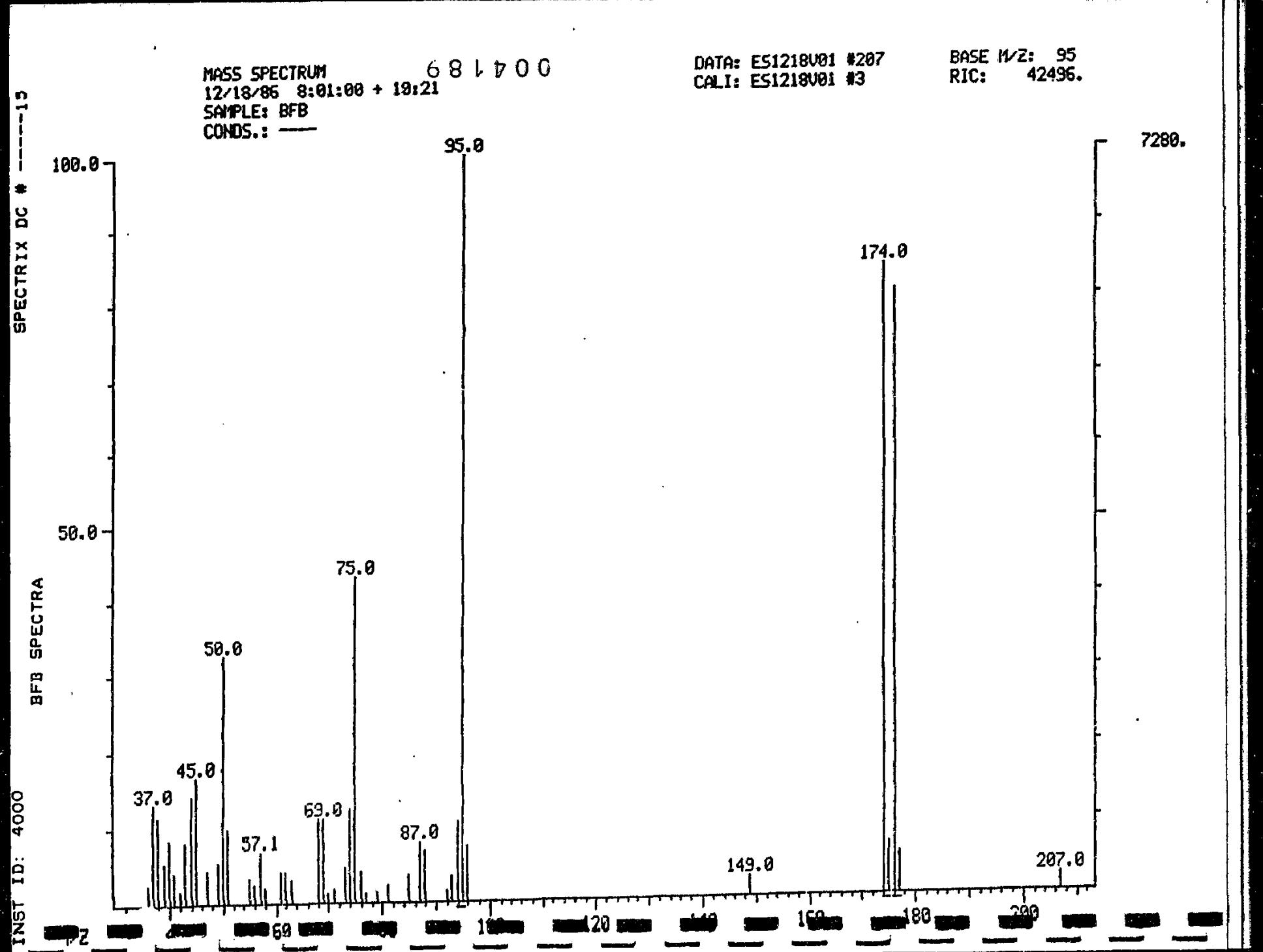
Calib: ES1218V01 # 3

Base m/z: 95

RIC: 42496.

	36 207 Mass	0.00 % RA	0.00 % RIC	0. #	Minima 0 Maxima	Min Inten: 0.
	35.97?	2.69	0.46	196.		
	36.99?	12.97	2.22	944.		
	37.98?	11.30	1.94	823.		
	39.01?	5.55	0.95	404.		
	39.95?	8.35	1.43	608.		
	41.02?	4.24	0.73	309.		
	42.02?	1.83	0.31	133.		
	43.04?	8.15	1.40	593.		
	43.99?	14.20	2.43	1034.		
	45.02?	16.51	2.83	1202.		
	47.02?	4.34	0.74	316.		
	49.01?	5.62	0.96	409.		
	50.01	32.73	5.61	2384.		
	51.02	9.57	1.64	697.		
	55.04	3.30	0.56	240.		
	56.04	2.71	0.46	197.		
	57.06	6.92	1.19	504.		
	58.07	2.17	0.37	158.		
	61.02	4.23	0.72	308.		
	62.05	4.07	0.70	296.		
	63.05	3.17	0.54	231.		
	58.02	10.88	1.86	792.		
	69.02	11.07	1.90	806.		
	70.05	1.40	0.24	102.		
	71.09	1.85	0.32	135.		
	73.02	4.60	0.79	335.		
	74.01	12.34	2.11	898.		
	75.02	43.08	7.38	3136.		
	76.03	4.08	0.70	297.		
	77.03	1.43	0.24	104.		
	78.91	1.52	0.26	111.		
	81.02	2.25	0.39	164.		
	84.99	3.72	0.64	271.		
	86.99	7.84	1.34	571.		
	87.97	6.84	1.17	498.		
	92.04	1.68	0.29	122.		
	93.02	3.39	0.58	247.		
	94.03 F	10.37	1.78	755.		
	95.02 F	100.00	17.13	7280.		
	96.02	7.23	1.24	528.		
	149.05	2.54	0.44	185.		
	173.95 F	84.40	14.46	6144.		
	174.95 F	6.91	1.18	503.		
	175.95 F	80.99	13.87	5896.		
	176.95 F	5.59	0.96	407.		
	207.03	2.49	0.43	181.		

0041100



GC/MS Tuning and Mass Calibration

BROMOFLUOROBENZENE

File Number:

Inst ID: 4000

Laboratory:

Sens Date: 12/19/86

Contract:

Lab ID: ES1219V01

Cali Date:

Sens Time: 6:48:00

Analyst: NOT

Data release authorized by:

m/z

Ion Abundance Criteria

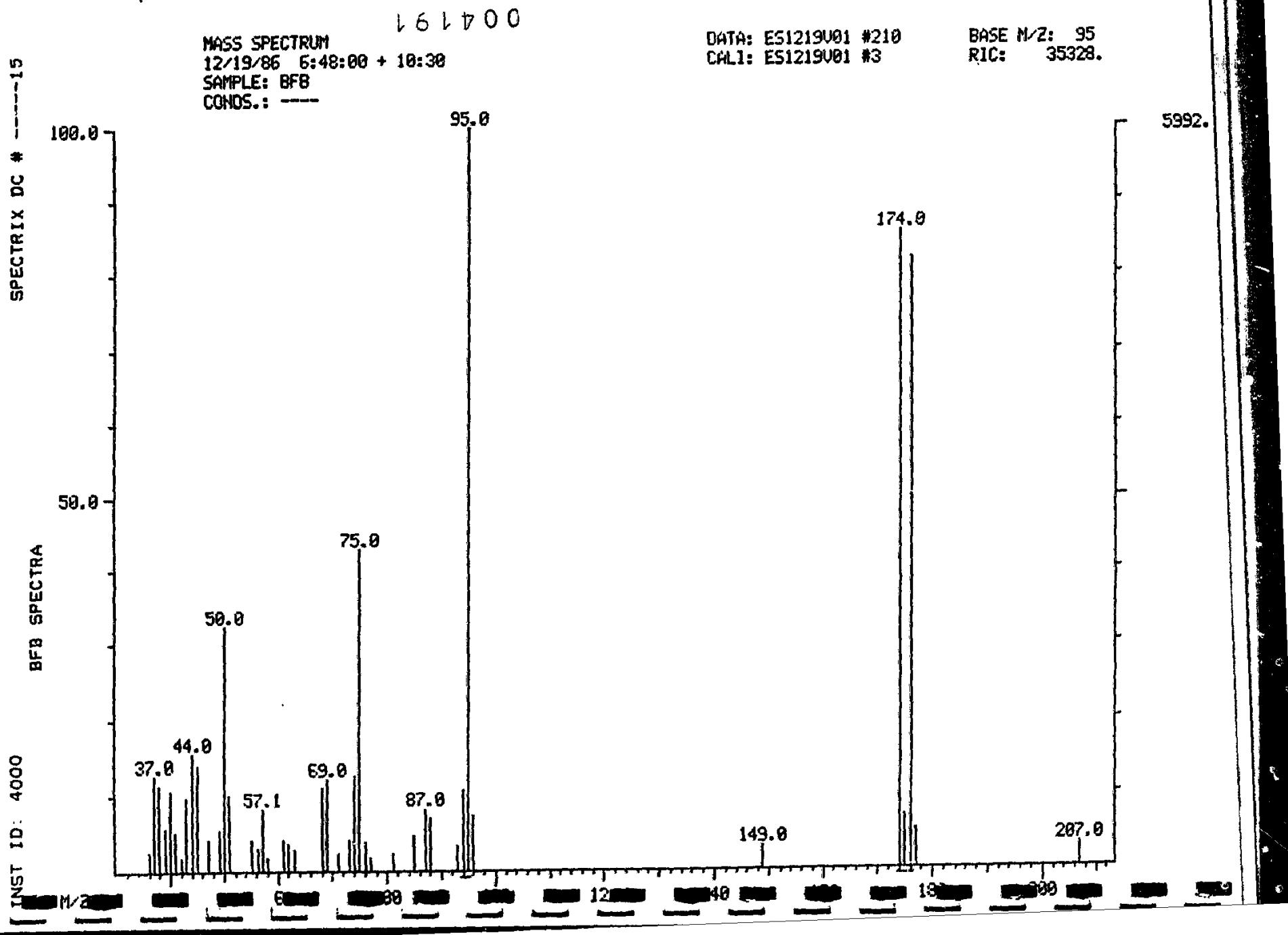
Spec # 210

50	15 to 40% of mass 95	32.51
75	30 to 60% of mass 95	42.99
95	base peak, 100% relative abundance	100.00
96	5 to 9% of mass 95	7.44
173	less than 1% of mass 95	0.00
174	greater than 50% of mass 95	83.85
175	5 to 9% of mass 174	7.09 (8.26) 1
176	between 95% and 101% of mass 174	82.11 (95.64) 1
177	5 to 9% of mass 176	5.31 (6.46) 2

1 - value in parenthesis is % of mass 174

2 - value in parenthesis is % of mass 176

004190



INST ID: 4000

BFB MASS LISTING

SPECTRIX DC # -----15

Mass List

12/19/86 6:48:00 + 10:30

Data: ES1219V01 # 210

Base m/z: 95

Sample: BFB

Calib: ES1219V01 # 3

RIC: 35328.

Rands.: ----

Mass	0.00	0.00	0. Minima # Inten.	Min Inten:	0.
	% RA	% RIC	0. Maxima		
35. 96?	2. 67	0. 45	160.		
36. 98?	12. 60	2. 14	795.		
37. 97?	11. 30	1. 92	677.		
39. 00?	5. 76	0. 98	345.		
39. 93?	10. 50	1. 78	629.		
41. 02?	5. 11	0. 87	306.		
42. 02?	1. 89	0. 32	113.		
43. 03?	9. 60	1. 63	575.		
43. 98	15. 49	2. 63	928.		
45. 01	13. 90	2. 36	833.		
47. 02	4. 14	0. 70	248.		
49. 01	5. 37	0. 91	322.		
50. 01	32. 51	5. 51	1948.		
51. 02	9. 86	1. 67	591.		
55. 05	4. 31	0. 73	258.		
56. 05	3. 07	0. 52	184.		
57. 07	8. 06	1. 37	483.		
58. 06	1. 92	0. 33	115.		
61. 03	4. 09	0. 69	245.		
62. 04	3. 69	0. 63	221.		
63. 05	3. 00	0. 51	180.		
68. 02	10. 98	1. 86	658.		
69. 02	11. 93	2. 02	715.		
71. 09	2. 39	0. 40	143.		
73. 03	4. 29	0. 73	257.		
74. 02	12. 47	2. 11	747.		
75. 02	42. 99	7. 29	2576.		
76. 05	3. 99	0. 68	239.		
77. 02	1. 80	0. 31	108.		
81. 02	2. 45	0. 42	147.		
84. 98	4. 72	0. 80	283.		
86. 98	8. 11	1. 38	486.		
87. 97	6. 94	1. 18	416.		
93. 01	3. 44	0. 58	206.		
94. 02 F	10. 76	1. 83	645.		
94. 99 F	100. 00	16. 96	5992.		
95. 98	7. 44	1. 26	446.		
149. 03	3. 04	0. 52	182.		
173. 95 F	85. 85	14. 56	5144.		
174. 97 F	7. 09	. 20	425.		
175. 95 F	82. 11	13. 93	4920.		
176. 97	5. 31	0. 90	318.		
207. 03	3. 09	0. 52	185.		

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